



No. 0000053

Gradient Corporation



**1989 Antrim Iron Works
(Tar Lake) Site Investigation**

Volume II - Tar and Soil Organic Compound Data

March 1990

44 Brattle Street □ Cambridge, MA 02138 □ (617) 576-1555 □ FAX: (617) 864-8469

EPA Region 5 Records Ctr.



209198

COMPUCHEM
LABORATORIES

JAN 19 1990

18/JAN/90

ENSAFE
ATTN. Mr. Phil Coop
P.O.Box341315(Zip38184-1315)
5705 Stage Rd./Suite 212
Memphis, TN 38134

Subject. Report of Data - Account Number 178202

ATTN. Mr. Phil Coop

Enclosed herewith are the results of analytical work and associated Quality Control data for the samples received by CompuChem@ under the referenced account number.

As you know, CLP service is performed in accordance with EPA methodologies as defined for the National Contract Laboratory Program; 'SOW For Organic Analysis; Multi-Media, Multi-Concentration' effective 2/88, revision 9/88. Your original data has been stored and is available for future reference.

Thank you for selecting CompuChem@ Laboratories for your sample analysis. We would like to continue providing you analytical support and services in the future. If you should have questions or require additional analytical services, please contact your representative at 1-919-549-8263.

Sincerely,



Elise L. Cobb
Supervisor, Report Deliverables

Attachment

COMPUCHEM
LABORATORIES, INC.

18/JAN/90

ENSAFE
ATTN. Mr. Phil Coop
P.O.Box341315(Zip38184-1315)
5705 Stage Rd./Suite 212
Memphis, TN 38134

ACCOUNT #. 178202

CC#	SAMPLE-ID	RECEIPT DATE
309679	B201A	12/20/89
309686	B201B	12/20/89
309687	B202TAR	12/20/89
309688	B202A	12/20/89
309689	B202B	12/20/89
309690	B202C	12/20/89

TOTAL NUMBER OF SAMPLES = 6

I. SAMPLE DATA SUMMARY PACKAGE

The Sample Data Summary Package shall contain data for samples in one Sample Delivery Group of the Case, as follows:

1. Case Narrative
2. By fraction (VOA, SV, PEST) and by sample within each fraction - tabulated target compound results (Form I) and tentatively identified compounds (Form I, TIC) (VOA and SV only)
3. By fraction (VOA, SV, PEST) - surrogate spike analysis results (Form II) by matrix (Water and/or Soil) and for soil, by concentration (Low or Medium)
4. By fraction (VOA, SV, PEST) - matrix spike / matrix spike duplicate results (Form III)
5. By fraction (VOA, SV, PEST) - blank data (Form IV) and tabulated results (Form I) including tentatively identified compounds (Form I, TIC) (VOA and SV only)
6. By fraction (VOA and SV only) - internal standard area data (Form VIII)

CASE#: 18756 SDG#: 07 SAS#: _____

1. Case Narrative

CLP CASE NARRATIVE--CASE # 18756
Contract No. Z-88 (REVS) SDG No. 07
CompuChem Laboratories, Inc.

Sample Numbers: B201A, B201B, B202A, B202B, B202C, B202TAR

This portion of Case # 18756 consisted of six solid samples for volatile and semivolatile analysis. The samples were received intact on 12/20/89 via Federal Express in properly sealed shipping containers with traffic reports. Due to a discrepancy between the sample container and the chain-of-custody, the C-O-C ID was used for sample B202A. Moisture content of the samples ranged from 9% to 64%. The pH values were within the values specified in EPA protocols.

VOLATILES:

Because of the nature of the samples, it was not possible to analyze the Low Level sample spikes within holding time requirements. All other volatile fractions were extracted and analyzed within holding time requirements. Because of the level of sample constituents, two analyses, a Low Level and a Medium Level, were reported for samples B201A, B201B, and B202TAR. The Low Level analyses of those three samples all contained toluene, 2-hexanone and total xylenes at levels outside their analytical ranges. In addition, sample B201A also contained benzene and ethylbenzene and sample B202TAR, methylene chloride, acetone, and ethylbenzene, at levels outside their analytical range in the Low Level analyses. Because of sample inhomogeneity and the differences in preparation procedures for the Low and Medium analyses, some of the reported concentrations did not compare well between the two reported analyses for these three samples. The total xylene concentration was mistakenly flagged with an "E" in the dilution of sample B202TAR. Since the analysis partially separates the xylenes, the individual xylene concentrations were acceptable. The surrogate DB-toluene failed recovery criteria in sample B201A due to matrix interference. All other surrogate recoveries met acceptance criteria. Both the Low and Medium Level matrix spike/matrix spike duplicate results were acceptable. Due to a scheduling error, the sample requested for use as the spike original was not used. Toluene failed recovery criteria in the Medium Level MS and MSD, and also failed %RPD in the Medium Level analyses.

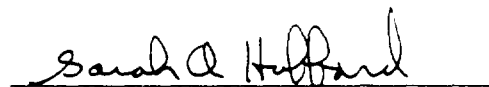
SEMIVOLATILES:

All semivolatile fractions were extracted and analyzed within holding time requirements. Coeluting compounds were flagged with an "X" on Form I. Because of the levels of 2-methylphenol, 4-methylphenol, and 2,4-dimethylphenol present, two analyses were reported for sample B201A, a 20:1 and a 50:1 dilution. Both a neat and a 5:1 analysis were reported for sample B202B because of the levels of 4-methylphenol and 2,4-dimethylphenol present. In addition to the above-mentioned TCL's, such compounds as phenol and various polyanomatic hydrocarbons were detected in one or more of these samples. Due to the viscosity of the extract, only a 5:1 dilution of the Medium Level extraction was reported for sample B201B. Because of the dilution necessary, no surrogate recovery data was available for the Low Level matrix spike/matrix spike duplicate. The recovery for OS-nitrobenzene in sample B201A exceeded recovery criteria. All other surrogates met recovery criteria. Because of the dilution necessary, there was no Low Level spike recovery data. A Low Level blank spike, extracted and analyzed with these samples, had acceptable results.

The failure of the %RPD for 2,4-dinitrotoluene in the blank spike was allowed, since the %RPD criteria are internal CompuChem requirements and not contractual necessities. The Medium Level matrix spike/matrix spike duplicate results were acceptable. The recoveries for n-nitroso-di-n-propylamine and 1,2,4-trichlorobenzene failed QC limits for both the MS and MSD. In addition, the recovery for 1,4-dichlorobenzene failed QC limits in the MSD. The %RPD's exceeded QC limits for 1,4-dichlorobenzene, 1,2,4-trichlorobenzene, and acenaphthene.

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or his designee as verified by the following signature.

Note: This report was paginated for reference and accountability in decreasing numerical sequence.



Sarah A. Hubbard 1/18/90
Technical Reviewer



COMPUCHEM
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DETECTION LIMIT CALCULATION CLARIFICATION

To protect our GC columns from unnecessary contamination, soil samples prepared according to Caucus Protocol methods are routinely diluted 5:1. Through a series of experiments we have determined that our Instrument Detection Limit for pesticides is 5X lower than the EPA Contract Required Quantitation Limit (CRQL). We, therefore, only adjust our detection limits if the dilution necessary to analyze the sample is greater than 5:1. If the sample is diluted by a factor of X the detection limit is adjusted by X/5 instead of X.

Doug M. McCormack

Doug McCormack
Manager, GC Laboratory

Robert E. Meierer

Robert E. Meierer
Vice - President, Quality Assurance



COMPUCHEM LABORATORIES

CHANGES TO THE EPA'S ORGANIC STATEMENT OF WORK (SOW) FOR THE CONTRACT LABORATORY PROGRAM (CLP)

Effective with samples received at CompuChem Laboratories, Inc. on April 3, 1989, the new "SOW for Organics Analysis; Multi-Media, Multi-Concentration" will be in effect. The new SOW is dated 2/88 with revisions dated 9/88.

EPA introduces modifications to the CLP SOW for a variety of reasons. They are:

- as a result of technical caucuses attended by representatives from EPA regional laboratories and the CLP, new or modified analytical methods are required,
- as a result of analytical data being supplied to the Agency by the CLP laboratories, QC acceptance criteria are updated and made a requirement of the program,
- as a result of requirements by the end users of the data (the EPA regions and the Program Office), changes to the deliverable requirements of the CLP are necessary.

As a service to our clients utilizing the EPA CLP SOW for their analytical needs, the following information is provided to point out the substantive changes between the new SOW and the previous one (10/86 with revisions through 8/87).

"KEY" CHANGES TO THE ORGANIC SOW

- 1) Wide bore capillary columns (internal diameter greater than 0.32 mm) are allowed for pesticide/PCB analysis in addition to packed columns. Pused silica capillary columns remain an optional confirmation column.

- 2) The Sample Data Summary Package (SDSP) was introduced in the 10/86 SOW. The SDSP consists of copies of specific summary forms. In the 2/88 SOW, Form VIII, the Internal Standard Area Summary Form for volatile and semi-volatile analyses, has been added to the SDSP. Form VIII for volatiles and semi-volatiles is also to be included in the QC Summary Package.
- 3) A clarification for dealing with the three xylene isomers has been added. For total xylenes, where three isomers are quantified as two peaks, the calibration range of each peak should be considered separately, e.g., a diluted analysis is not required for total xylenes unless the concentration of either peak exceeds 200 ug/l.
- 4) The volatile target analyte, 2-butanone, is to be quantified against the first internal standard, bromochloromethane. In the prior SOW, 2-butanone was quantified against the second internal standard, 1,4-difluorobenzene.
- 5) Turnaround time for the delivery of data has been lowered from 40 days to 35 days for routine analytical service (RAS) work associated with the Superfund program. The turnaround time is calculated from the Validated Time of Sample Receipt (VTSR) of the last sample in the Sample Delivery Group (SDG).

The above represents the major changes in the new organic SOW for EPA's CLP. There are, of course, other minor changes which have not been addressed in this announcement. Of note, the hardcopy deliverable forms have not changed and, in fact, retain the same revision dates as the prior SOW.

If there are any questions concerning the information presented, please feel free to contact your account administrator at 1-800-833-8097.



Robert E. Meierer
Vice President of Quality Assurance

DATA REPORTING QUALIFIERS

For reporting results to EPA, the following result qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

VALUE - If the result is a value greater than or equal to the detection limit, report the value.

U - Indicates compound was analyzed but not detected. The sample quantitation limit must be corrected for dilution and for percent moisture. For example, 10 U for phenol in water if the sample final volume is the protocol-specified final volume. If a 1 to 10 dilution of extract is necessary, the reported limit is 100 U. For a soil sample, the value must also be adjusted for percent moisture. For example, if the sample had 24% moisture and a 1 to 10 dilution factor, the sample quantitation limit for phenol (330 U) would be corrected to:

$$\frac{(330 \text{ U}) \times df}{D} \quad \text{where } D = \frac{100 - \% \text{ moisture}}{100}$$

and df = dilution factor

$$\text{at } 24\% \text{ moisture, } D = \frac{100-24}{100} = 0.76$$

$$\frac{(330 \text{ U}) \times 10}{.76} = 4300 \text{ U} \quad \text{rounded to the appropriate number of significant figures}$$

For soil sample subjected to BPC clean-up procedures, the CRDL is also multiplied by 2, to account for the fact that only half of the extract is recovered.

J - Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed, or when the mass spectral data indicate the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than zero. For example, if the sample quantitation limit is 10 ug/L, but a concentration of 3 ug/L is calculated, report it as 3J. The sample quantitation limit must be adjusted for both dilution and percent moisture as discussed for the U flag, so that if a sample with 24% moisture and a 1 to 10 dilution factor has a calculated concentration of 300 ug/L and a sample quantitation limit of 430 ug/kg, report the concentration as 300J on Form 1.

C - This flag applies to pesticide results where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ug/l in the final extract shall be confirmed by GC/MS.

DATA REPORTING QUALIFIERS - PAGE 2

- B** - This flag is used when the analyte is found in the associated blank as well as in the sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action. This flag must be used for a TIC as well as for a positively identified TCL compound.

- E** - This flag identifies compounds whose concentrations exceed the calibration range of the GC/MS instrument for that specific analysis. This flag will not apply to pesticides/PCBs analyzed by GC/EC methods. If one or more compounds have a response greater than full scale, the sample or extract must be diluted and re-analyzed according to the specifications. All such compounds with a response greater than full scale should have the concentration flagged with an "E" on the Form I for the original analysis. If the dilution of the extract causes any compounds identified in the first analysis to be below the calibration range in the second analysis, then the results of both analyses shall be reported on separate Forms I. The Form I for the diluted sample shall have the "DL" suffix appended to the sample number.

- D** - This flag identifies all compounds identified in an analysis at a secondary dilution factor. If a sample or extract is re-analyzed at a higher dilution factor, as in the "E" flag above, the "DL" suffix is appended to the sample number on the Form I for the diluted sample and all concentration values reported on that Form I are flagged with the "D" flag.

- A** - This flag indicates that TIC is a suspected aldol-condensation product.

- X** - Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the Sample Data Summary Package and the Case Narrative. If more than one is required, use "Y" and "Z", as needed. If more than five qualifiers are required for a sample result, use the "X" flag to combine several flags, as needed. For instance, the "X" flag might combine the "A", "B", and "D" flags for some sample.




COMPUCHEM
LABORATORIES

QUALITY ASSURANCE NOTICE

With the advent of the new organics Statement of Work (SOW 2/88, Revision: 9/88) participants in EPA's Contract Laboratory Program (CLP) are required to provide hard copy and diskette deliverables. CompuChem employs the Finnigan QA Formaster Program (Format A) to generate these requirements using data files from our analytical instrumentation. Currently, and independently, quantitation reports are generated by the instruments and are used with CompuChem-developed software to calculate results. The GC and GC/MS quantitation routines employ the convention of carrying at least one extra significant figure until the mathematical computations are completed. Then, the quantitative results are rounded to the SOW-required number of significant figures for reporting. In addition, the algorithm used by the Formaster Program is slightly different than that employed in CompuChem's software routines. Therefore, results presented in the supportive data supplied with our deliverables packages may be slightly different than those which appear on the hard copy forms generated via Formaster.

This notice serves to alert the end users of these data packages as to the reason why slight differences may be observed.



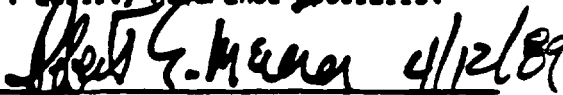
Robert E. Meierer
Director of Quality Assurance

QUALITY ASSURANCE NOTICE

Specific guidelines are presented in the EPA CLP Organic Statement of Work for the positive qualitative identification of compounds through mass spectral interpretation. Applying these guidelines absolutely may not be possible when the nature of the sample is less than pure reference material. Where the mass spectral pattern of a compound to be identified demonstrates interferences or resolution from one or more additional compounds, either unknowns, internal standards, or surrogate standards, the "+" sign is added to the top of the dual spectra page.

 4/10/89

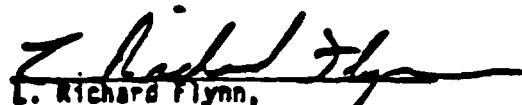
Linda Fowler
Sr. Quality Assurance Specialist

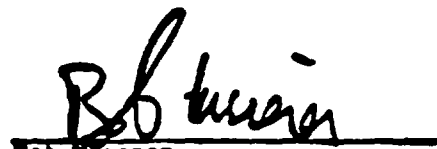
 4/12/89

Robert E. Malerer
Vice President, Quality Assurance

LABORATORY NOTICE

On June 15, 1985 CompuChem Laboratories began adding D3-2,4-Dinitrophenol to all standards and samples. The purpose of this addition is to enable the laboratory to have higher and more consistent analytical sensitivity for the native 2,4-Dinitrophenol. The peak corresponding to the deuterated analog is clearly labeled on each RIC as D3/1 and will not be searched and reported as a tentatively identified compound (TIC). This compound is not being used as an internal or surrogate standard.


L. Richard Flynn,
Development Chemist


Bob Krierer,
Director of Quality Assurance

2. By fraction (VOA , SV , PEST) and by sample within each fraction - tabulated target compound results (Form I) and tentatively identified compounds (Form I, TIC) (VOA and SV only)

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: 309679

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009679C19

Level: (low/med) LOW Date Received: 12/20/89

% Moisture: not dec. 9 Date Analyzed: 12/22/89

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	11	U
74-83-9	-----Bromomethane	11	U
75-01-4	-----Vinyl Chloride	11	U
75-00-3	-----Chloroethane	11	U
75-09-2	-----Methylene Chloride	32	B
67-64-1	-----Acetone	150	B
75-15-0	-----Carbon Disulfide	5	U
75-35-4	-----1,1-Dichloroethene	5	U
75-34-3	-----1,1-Dichloroethane	5	U
540-59-0	-----1,2-Dichloroethene (total)	5	U
67-66-3	-----Chloroform	5	U
107-06-2	-----1,2-Dichloroethane	5	U
78-93-3	-----2-Butanone	120	U
71-55-6	-----1,1,1-Trichloroethane	5	U
56-23-5	-----Carbon Tetrachloride	5	U
108-05-4	-----Vinyl Acetate	11	U
75-27-4	-----Bromodichloromethane	5	U
78-87-5	-----1,2-Dichloropropane	5	U
10061-01-5	-----cis-1,3-Dichloropropene	5	U
79-01-6	-----Trichloroethene	5	U
124-48-1	-----Dibromochloromethane	5	U
79-00-5	-----1,1,2-Trichloroethane	5	U
71-43-2	-----Benzene	260	E
10061-02-6	-----Trans-1,3-Dichloropropene	5	U
75-25-2	-----Bromoform	5	U
108-10-1	-----4-Methyl-2-Pentanone	17	U
591-78-6	-----2-Hexanone	580	E
127-18-4	-----Tetrachloroethene	5	U
79-34-5	-----1,1,2,2-Tetrachloroethane	5	U
108-88-3	-----Toluene	720	E
108-90-7	-----Chlorobenzene	5	U
100-41-4	-----Ethylbenzene	420	E
100-42-5	-----Styrene	210	U
1330-20-7	-----Total Xylenes	1300	E

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009679C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 625-86-5	FURAN, 2,5-DIMETHYL-	8.22	240	J
2. 1192-62-7	ETHANONE, 1-(2-FURANYL)-	11.02	250	J
3.	UNKNOWN	13.44	460	J
4.	UNKNOWN	13.45	220	J
5.	UNKNOWN	14.05	190	J
6.	UNKNOWN	14.55	300	J
7. 100-66-3	BENZENE, METHOXY-	14.70	510	J
8. 103-65-1	BENZENE, PROPYL-	15.34	190	J
9.	UNKNOWN	15.60	1500	J
10.	SUBSTITUTED BENZENE	16.05	800	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201ARE

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C3R09679B13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 9 Date Analyzed: 12/28/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	Chloromethane	1400	U
74-83-9	Bromomethane	1400	U
75-01-4	Vinyl Chloride	1400	U
75-00-3	Chloroethane	1400	U
75-09-2	Methylene Chloride	440	BDJ
67-64-1	Acetone	1400	U
75-15-0	Carbon Disulfide	690	U
75-35-4	1,1-Dichloroethene	690	U
75-34-3	1,1-Dichloroethane	690	U
540-59-0	1,2-Dichloroethene (total)	690	U
67-66-3	Chloroform	690	U
107-06-2	1,2-Dichloroethane	690	U
78-93-3	2-Butanone	1400	U
71-55-6	1,1,1-Trichloroethane	690	U
56-23-5	Carbon Tetrachloride	690	U
108-05-4	Vinyl Acetate	1400	U
75-27-4	Bromodichloromethane	690	U
78-87-5	1,2-Dichloropropane	690	U
10061-01-5	cis-1,3-Dichloropropene	690	U
79-01-6	Trichloroethene	690	U
124-48-1	Dibromochloromethane	690	U
79-00-5	1,1,2-Trichloroethane	690	U
71-43-2	Benzene	630	DJ
10061-02-6	Trans-1,3-Dichloropropene	690	U
75-25-2	Bromoform	690	U
108-10-1	4-Methyl-2-Pentanone	1400	U
591-78-6	2-Hexanone	630	DJ
127-18-4	Tetrachloroethene	690	U
79-34-5	1,1,2,2-Tetrachloroethane	690	U
108-88-3	Toluene	4200	D
108-90-7	Chlorobenzene	690	U
100-41-4	Ethylbenzene	3100	D
100-42-5	Styrene	1300	D
1330-20-7	Total Xylenes	9500	D

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201ARE

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C3R09679B13
 Level: (low/med; MED) Date Received: 12/20/89
 % Moisture: not dec. 9 Date Analyzed: 12/28/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 9 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 1192-62-7	ETHANONE, 1-(2-FURANYL) -	9.80	820	J
2.	UNKNOWN	12.19	1200	J
3.	SUBSTITUTED BENZENE	14.05	1400	J
4.	SUBSTITUTED BENZENE	14.29	8400	J
5.	SUBSTITUTED BENZENE	14.45	2300	J
6.	SUBSTITUTED BENZENE	14.77	7000	J
7.	SUBSTITUTED BENZENE	15.14	9900	J
8. 496-11-7	1H-INDENE, 2, 3-DIHYDRO-	15.30	960	J
9. 271-89-6	BENZOFURAN	15.55	9200	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009686C19
 Level: (low/med) LOW Date Received: 12/20/89
 ‡ Moisture: not dec. 12 Date Analyzed: 12/27/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	57	U
74-83-9	-----Bromomethane	57	U
75-01-4	-----Vinyl Chloride	57	U
75-00-3	-----Chloroethane	57	U
75-09-2	-----Methylene Chloride	69	B
67-64-1	-----Acetone	370	B
75-15-0	-----Carbon Disulfide	28	U
75-35-4	-----1,1-Dichloroethene	28	U
75-34-3	-----1,1-Dichloroethane	28	U
540-59-0	-----1,2-Dichloroethene (total)	28	U
67-66-3	-----Chloroform	28	U
107-06-2	-----1,2-Dichloroethane	28	U
78-93-3	-----2-Butanone	260	
71-55-6	-----1,1,1-Trichloroethane	28	U
56-23-5	-----Carbon Tetrachloride	28	U
108-05-4	-----Vinyl Acetate	57	U
75-27-4	-----Bromodichloromethane	28	U
78-87-5	-----1,2-Dichloropropane	28	U
10061-01-5	-----cis-1,3-Dichloropropene	28	U
79-01-6	-----Trichloroethene	28	U
124-48-1	-----Dibromochloromethane	28	U
79-00-5	-----1,1,2-Trichloroethane	28	U
71-43-2	-----Benzene	140	
10061-02-6	-----Trans-1,3-Dichloropropene	28	U
75-25-2	-----Bromoform	28	U
108-10-1	-----4-Methyl-2-Pentanone	71	
591-78-6	-----2-Hexanone	1600	BE
127-18-4	-----Tetrachloroethene	28	U
79-34-5	-----1,1,2,2-Tetrachloroethane	28	U
108-88-3	-----Toluene	1600	E
108-90-7	-----Chlorobenzene	28	U
100-41-4	-----Ethylbenzene	980	
100-42-5	-----Styrene	540	
1330-20-7	-----Total Xylenes	3300	E

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1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009686C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/27/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	11.04	500	J
2.	METHYLESTERPENTENOIC ACID +	13.45	1400	J
3.	UNKNOWN	14.09	1100	J
4. 110-43-0	2-HEPTANONE	14.27	1400	J
5.	UNKNOWN	14.42	910	J
6.	UNKNOWN	14.55	1100	J
7. 100-66-3	BENZENE, METHOXY-	14.72	2300	J
8.	UNKNOWN	15.35	500	J
9.	ETHYLMETHYLBENZENE + UNKNOWN	15.60	5400	J
10. 611-14-3	BENZENE, 1-ETHYL-2-METHYL-	16.07	1500	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201BRE

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C5R09686B13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/29/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	1400	U
74-83-9	-----Bromomethane	1400	U
75-01-4	-----Vinyl Chloride	1400	U
75-00-3	-----Chloroethane	1400	U
75-09-2	-----Methylene Chloride	540	DBJ
67-64-1	-----Acetone	1400	U
75-15-0	-----Carbon Disulfide	710	U
75-35-4	-----1,1-Dichloroethene	710	U
75-34-3	-----1,1-Dichloroethane	710	U
540-59-0	-----1,2-Dichloroethene (total)	710	U
67-66-3	-----Chloroform	710	U
107-06-2	-----1,2-Dichloroethane	710	U
78-93-3	-----2-Butanone	1400	U
71-55-6	-----1,1,1-Trichloroethane	710	U
56-23-5	-----Carbon Tetrachloride	710	U
108-05-4	-----Vinyl Acetate	1400	U
75-27-4	-----Bromodichloromethane	710	U
78-87-5	-----1,2-Dichloropropane	710	U
10061-01-5	-----cis-1,3-Dichloropropene	710	U
79-01-6	-----Trichloroethene	710	U
124-48-1	-----Dibromochloromethane	710	U
79-00-5	-----1,1,2-Trichloroethane	710	U
71-43-2	-----Benzene	2200	D
10061-02-6	-----Trans-1,3-Dichloropropene	710	U
75-25-2	-----Bromoform	710	U
108-10-1	-----4-Methyl-2-Pentanone	1400	U
591-78-6	-----2-Hexanone	2400	D
127-18-4	-----Tetrachloroethene	710	U
79-34-5	-----1,1,2,2-Tetrachloroethane	710	U
108-88-3	-----Toluene	16000	D
108-90-7	-----Chlorobenzene	710	U
100-41-4	-----Ethylbenzene	7400	D
100-42-5	-----Styrene	2800	D
1330-20-7	-----Total Xylenes	21000	D

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VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201BRE

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C5R09686B13
 Level: (low/med; MED Date Received: 12/20/89
 ‡ Moisture: not dec. 12 Date Analyzed: 12/29/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 625-86-5	FURAN, 2, 5-DIMETHYL-	7.28	1800	J
2.	UNKNOWN	10.02	2000	J
3.	UNKNOWN	12.34	2800	J
4.	METHYLETHYLBENZENE + UNKNOWN	13.44	710	J
5. 103-65-1	BENZENE, PROPYL-	14.22	3100	J
6. 98-82-8	BENZENE, (1-METHYLETHYL) -	14.45	17000	J
7.	TRIMETHYLBENZENE ISOMER	14.60	4700	J
8.	ETHYLMETHYLBENZENE + UNKNOWN	14.94	14000	J
9.	TRIMETHYLBENZENE ISOMER	15.30	18000	J
10. 611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	15.45	2000	J

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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309688
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009688C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/22/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	Chloromethane	11	U
74-83-9	Bromomethane	11	U
75-01-4	Vinyl Chloride	11	U
75-00-3	Chloroethane	11	U
75-09-2	Methylene Chloride	33	B
67-64-1	Acetone	37	B
75-15-0	Carbon Disulfide	6	U
75-35-4	1,1-Dichloroethene	6	U
75-34-3	1,1-Dichloroethane	6	U
540-59-0	1,2-Dichloroethene (total)	6	U
67-66-3	Chloroform	6	U
107-06-2	1,2-Dichloroethane	6	U
78-93-3	2-Butanone	26	
71-55-6	1,1,1-Trichloroethane	6	U
56-23-5	Carbon Tetrachloride	6	U
108-05-4	Vinyl Acetate	11	U
75-27-4	Bromodichloromethane	6	U
78-87-5	1,2-Dichloropropane	6	U
10061-01-5	cis-1,3-Dichloropropene	6	U
79-01-6	Trichloroethene	6	U
124-48-1	Dibromochloromethane	6	U
79-00-5	1,1,2-Trichloroethane	6	U
71-43-2	Benzene	6	U
10061-02-6	Trans-1,3-Dichloropropene	6	U
75-25-2	Bromoform	6	U
108-10-1	4-Methyl-2-Pentanone	11	U
591-78-6	2-Hexanone	13	
127-18-4	Tetrachloroethene	6	U
79-34-5	1,1,2,2-Tetrachloroethane	6	U
108-88-3	Toluene	1	J
108-90-7	Chlorobenzene	6	U
100-41-4	Ethylbenzene	6	U
100-42-5	Styrene	6	U
1330-20-7	Total Xylenes	2	J

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VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309688
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009688C19
 Level: (low/med; LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 4

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	13.00	6.8	J
2.	UNKNOWN	14.37	5.7	J
3.	UNKNOWN	15.57	6.8	J
4.	INSTRUMENT ARTIFACT	15.92	6.8	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009689C19
 Level: (low/mc³) LOW Date Received: 12/20/89
 % Moisture: not dec. 14 Date Analyzed: 12/22/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	12	U
74-83-9	-----Bromomethane	12	U
75-01-4	-----Vinyl Chloride	12	U
75-00-3	-----Chloroethane	12	U
75-09-2	-----Methylene Chloride	30	B
67-64-1	-----Acetone	53	B
75-15-0	-----Carbon Disulfide	6	U
75-35-4	-----1,1-Dichloroethene	6	U
75-34-3	-----1,1-Dichloroethane	6	U
540-59-0	-----1,2-Dichloroethene (total)	6	U
67-66-3	-----Chloroform	6	U
107-06-2	-----1,2-Dichloroethane	6	U
78-93-3	-----2-Butanone	40	
71-55-6	-----1,1,1-Trichloroethane	6	U
56-23-5	-----Carbon Tetrachloride	6	U
108-05-4	-----Vinyl Acetate	12	U
75-27-4	-----Bromodichloromethane	6	U
78-87-5	-----1,2-Dichloropropane	6	U
10061-01-5	-----cis-1,3-Dichloropropene	6	U
79-01-6	-----Trichloroethene	6	U
124-48-1	-----Dibromochloromethane	6	U
79-00-5	-----1,1,2-Trichloroethane	6	U
71-43-2	-----Benzene	6	U
10061-02-6	-----Trans-1,3-Dichloropropene	6	U
75-25-2	-----Bromoform	6	U
108-10-1	-----4-Methyl-2-Pentanone	12	U
591-78-6	-----2-Hexanone	23	
127-18-4	-----Tetrachloroethene	6	U
79-34-5	-----1,1,2,2-Tetrachloroethane	6	U
108-88-3	-----Toluene	15	
108-90-7	-----Chlorobenzene	6	U
100-41-4	-----Ethylbenzene	12	
100-42-5	-----Styrene	3	J
1330-20-7	-----Total Xylenes	54	

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VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009689C19
 Level: (low/mcd) LOW Date Received: 12/20/89
 % Moisture: not dec. 14 Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 7

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	13.39	8.1	J
2.	UNKNOWN	14.05	7.0	J
3.	UNKNOWN	14.40	8.1	J
4.	UNKNOWN	14.49	8.1	J
5. 103-65-1	BENZENE, PROPYL-	15.35	8.1	J
6. 620-14-4	BENZENE, 1-ETHYL-3-METHYL-	15.55	84	J
7. 611-14-3	BENZENE, 1-ETHYL-2-METHYL-	16.04	21	J

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EPA SAMPLE NO.

B202C

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309690
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009690A19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 13 Date Analyzed: 12/22/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	11	U
74-83-9	-----Bromomethane	11	U
75-01-4	-----Vinyl Chloride	11	U
75-00-3	-----Chloroethane	11	U
75-09-2	-----Methylene Chloride	23	B
67-64-1	-----Acetone	26	B
75-15-0	-----Carbon Disulfide	6	U
75-35-4	-----1,1-Dichloroethene	6	U
75-34-3	-----1,1-Dichloroethane	6	U
540-59-0	-----1,2-Dichloroethene (total)	6	U
67-66-3	-----Chloroform	6	U
107-06-2	-----1,2-Dichloroethane	6	U
78-93-3	-----2-Butanone	3	J
71-55-6	-----1,1,1-Trichloroethane	6	U
56-23-5	-----Carbon Tetrachloride	6	U
108-05-4	-----Vinyl Acetate	11	U
75-27-4	-----Bromodichloromethane	6	U
78-87-5	-----1,2-Dichloropropane	6	U
10061-01-5	-----cis-1,3-Dichloropropene	6	U
79-01-6	-----Trichloroethene	6	U
124-48-1	-----Dibromochloromethane	6	U
79-00-5	-----1,1,2-Trichloroethane	6	U
71-43-2	-----Benzene	6	U
10061-02-6	-----Trans-1,3-Dichloropropene	6	U
75-25-2	-----Bromoform	6	U
108-10-1	-----4-Methyl-2-Pentanone	11	U
591-78-6	-----2-Hexanone	11	U
127-18-4	-----Tetrachloroethene	6	U
79-34-5	-----1,1,2,2-Tetrachloroethane	6	U
108-88-3	-----Toluene	2	J
108-90-7	-----Chlorobenzene	6	U
100-41-4	-----Ethylbenzene	1	J
100-42-5	-----Styrene	6	U
1330-20-7	-----Total Xylenes	7	

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VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202C

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309690
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009690A19
 Level: (low/med, LOW) Date Received: 12/20/89
 % Moisture: not dec. 13 Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 3 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	13.02	9.2	J
2. 4971-18-0	CYCLOPENTANONE, 2-ETHYL-	15.59	14	J
3.	INSTRUMENT ARTIFACT	15.97	9.2	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TAR

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: 309687

Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009687C19

Level: (low/med) LOW Date Received: 12/20/89

% Moisture: not dec. 64 Date Analyzed: 12/27/89

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	Chloromethane	140	U
74-83-9	Bromomethane	140	U
75-01-4	Vinyl Chloride	140	U
75-00-3	Chloroethane	140	U
75-09-2	Methylene Chloride	8300	EB
67-64-1	Acetone	3500	EB
75-15-0	Carbon Disulfide	69	U
75-35-4	1,1-Dichloroethene	69	U
75-34-3	1,1-Dichloroethane	69	U
540-59-0	1,2-Dichloroethene (total)	69	U
67-66-3	Chloroform	69	U
107-06-2	1,2-Dichloroethane	69	U
78-93-3	2-Butanone	5000	E
71-55-6	1,1,1-Trichloroethane	69	U
56-23-5	Carbon Tetrachloride	69	U
108-05-4	Vinyl Acetate	140	U
75-27-4	Bromodichloromethane	69	U
78-87-5	1,2-Dichloropropane	69	U
10061-01-5	cis-1,3-Dichloropropene	69	U
79-01-6	Trichloroethene	69	U
124-48-1	Dibromochloromethane	69	U
79-00-5	1,1,2-Trichloroethane	69	U
71-43-2	Benzene	1200	U
10061-02-6	Trans-1,3-Dichloropropene	69	U
75-25-2	Bromoform	69	U
108-10-1	4-Methyl-2-Pentanone	1200	U
591-78-6	2-Hexanone	12000	EB
127-18-4	Tetrachloroethene	69	U
79-34-5	1,1,2,2-Tetrachloroethane	69	U
108-88-3	Toluene	8600	E
108-90-7	Chlorobenzene	69	U
100-41-4	Ethylbenzene	8000	E
100-42-5	Styrene	2300	U
1330-20-7	Total Xylenes	15000	E

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TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202TAR

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009687C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 64 Date Analyzed: 12/27/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 534-22-5	FURAN, 2-METHYL-	4.95	9000	J
2. 625-86-5	FURAN, 2,5-DIMETHYL-	8.23	13000	J
3.	UNKNOWN	11.05	4700	J
4.	UNKNOWN	11.50	3200	J
5.	UNKNOWN	13.49	2600	J
6.	UNKNOWN	14.09	2900	J
7. 13368-65-5	CYCLOHEXANONE, 3-METHYL-, (R)-	14.45	1700	J
8.	UNKNOWN	14.75	3900	J
9.	UNKNOWN	15.65	11000	J
10. 3141-02-4	1,3-CYCLOPENTADIENE, 5-(1-MET	16.10	4400	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TARDL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C3R09687C13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 Date Analyzed: 12/29/89
 Column: (pack/cap) CAP Dilution Factor: 3.3

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	11000	U
74-83-9	-----Bromomethane	11000	U
75-01-4	-----Vinyl Chloride	11000	U
75-00-3	-----Chloroethane	11000	U
75-09-2	-----Methylene Chloride	16000	BD
67-64-1	-----Acetone	15000	D
75-15-0	-----Carbon Disulfide	5700	U
75-35-4	-----1,1-Dichloroethene	5700	U
75-34-3	-----1,1-Dichloroethane	5700	U
540-59-0	-----1,2-Dichloroethene (total)	5700	U
67-66-3	-----Chloroform	5700	U
107-06-2	-----1,2-Dichloroethane	5700	U
78-93-3	-----2-Butanone	11000	U
71-55-6	-----1,1,1-Trichloroethane	5700	U
56-23-5	-----Carbon Tetrachloride	5700	U
108-05-4	-----Vinyl Acetate	11000	U
75-27-4	-----Bromodichloromethane	5700	U
78-87-5	-----1,2-Dichloropropane	5700	U
10061-01-5	-----cis-1,3-Dichloropropene	5700	U
79-01-6	-----Trichloroethene	5700	U
124-48-1	-----Dibromochloromethane	5700	U
79-00-5	-----1,1,2-Trichloroethane	5700	U
71-43-2	-----Benzene	5000	DJ
10061-02-6	-----Trans-1,3-Dichloropropene	5700	U
75-25-2	-----Bromoform	5700	U
108-10-1	-----4-Methyl-2-Pentanone	11000	U
591-78-6	-----2-Hexanone	11000	U
127-18-4	-----Tetrachloroethene	5700	U
79-34-5	-----1,1,2,2-Tetrachloroethane	5700	U
108-88-3	-----Toluene	100000	D
108-90-7	-----Chlorobenzene	5700	U
100-41-4	-----Ethylbenzene	100000	D
100-42-5	-----Styrene	22000	D
1330-20-7	-----Total Xylenes	280000	DE

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VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202TARDL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C3R09687C13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 Date Analyzed: 12/29/89
 Column (pack/cap) CAP Dilution Factor: 3.3

Number TICs found: 10 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 534-22-5	FURAN, 2-METHYL-	4.20	29000	J
2. 625-86-5	FURAN, 2, 5-DIMETHYL-	7.15	49000	J
3. 1192-62-7	ETHANONE, 1-(2-FURANYL) -	9.84	61000	J
4. 37609-41-9	BICYCLO[3.2.1]OCT-2-ENE, 3-(1	12.24	85000	J
5. 24156-95-4	2-CYCLOPENTEN-1-ONE, 3, 5, 5-TR	12.80	40000	J
6.	UNKNOWN	13.30	34000	J
7.	SUBST. BENZENE	14.10	48000	J
8.	SUBST. BENZENE	14.35	300000	J
9.	SUBST. BENZENE	14.82	220000	J
10.	SUBST. BENZENE	15.19	170000	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202AMS

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309680
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: G3R09680A19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 01/02/90
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>		Q
74-87-3	-----Chloromethane	11		U
74-83-9	-----Bromomethane	11		U
75-01-4	-----Vinyl Chloride	11		U
75-00-3	-----Chloroethane	11		U
75-09-2	-----Methylene Chloride	16		B
67-64-1	-----Acetone	37		
75-15-0	-----Carbon Disulfide	6		U
75-35-4	-----1,1-Dichloroethene	6		U
75-34-3	-----1,1-Dichloroethane	6		U
540-59-0	-----1,2-Dichloroethene (total)	6		U
67-66-3	-----Chloroform	6		U
107-06-2	-----1,2-Dichloroethane	6		U
78-93-3	-----2-Butanone	20		
71-55-6	-----1,1,1-Trichloroethane	6		U
56-23-5	-----Carbon Tetrachloride	6		U
108-05-4	-----Vinyl Acetate	11		U
75-27-4	-----Bromodichloromethane	6		U
78-87-5	-----1,2-Dichloropropane	6		U
10061-01-5	-----cis-1,3-Dichloropropene	6		U
79-01-6	-----Trichloroethene	6		U
124-48-1	-----Dibromochloromethane	6		U
79-00-5	-----1,1,2-Trichloroethane	6		U
71-43-2	-----Benzene	6		U
10061-02-6	-----Trans-1,3-Dichloropropene	6		U
75-25-2	-----Bromoform	6		U
108-10-1	-----4-Methyl-2-Pentanone	11		U
591-78-6	-----2-Hexanone	8		J
127-18-4	-----Tetrachloroethene	6		U
79-34-5	-----1,1,2,2-Tetrachloroethane	6		U
108-88-3	-----Toluene	6		U
108-90-7	-----Chlorobenzene	6		U
100-41-4	-----Ethylbenzene	6		U
100-42-5	-----Styrene	6		U
1330-20-7	-----Total Xylenes	6		U

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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201BREMS

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 311271
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: CN011271B13
 Level: (low/med) MED Date Received: 12/20/89
 ‡ Moisture: not dec. 12 Date Analyzed: 12/29/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	1400	U
74-83-9	-----Bromomethane	1400	U
75-01-4	-----Vinyl Chloride	1400	U
75-00-3	-----Chloroethane	1400	U
75-09-2	-----Methylene Chloride	490	BJ
67-64-1	-----Acetone	1400	U
75-15-0	-----Carbon Disulfide	710	U
75-35-4	-----1,1-Dichloroethene	710	U
75-34-3	-----1,1-Dichloroethane	710	U
540-59-0	-----1,2-Dichloroethene (total)	710	U
67-66-3	-----Chloroform	710	U
107-06-2	-----1,2-Dichloroethane	710	U
78-93-3	-----2-Butanone	1400	U
71-55-6	-----1,1,1-Trichloroethane	710	U
56-23-5	-----Carbon Tetrachloride	710	U
108-05-4	-----Vinyl Acetate	1400	U
75-27-4	-----Bromodichloromethane	710	U
78-87-5	-----1,2-Dichloropropane	710	U
10061-01-5	-----cis-1,3-Dichloropropene	710	U
79-01-6	-----Trichloroethene	710	U
124-48-1	-----Dibromochloromethane	710	U
79-00-5	-----1,1,2-Trichloroethane	710	U
71-43-2	-----Benzene	710	U
10061-02-6	-----Trans-1,3-Dichloropropene	710	U
75-25-2	-----Bromoform	710	U
108-10-1	-----4-Methyl-2-Pentanone	1400	U
591-78-6	-----2-Hexanone	2700	
127-18-4	-----Tetrachloroethene	710	U
79-34-5	-----1,1,2,2-Tetrachloroethane	710	U
108-88-3	-----Toluene	710	U
108-90-7	-----Chlorobenzene	710	U
100-41-4	-----Ethylbenzene	7600	
100-42-5	-----Styrene	2900	
1330-20-7	-----Total Xylenes	22000	

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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202AMSD

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309681
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: G3R09681A19
 Level: (low/med) LOW Date Received: 12/20/89
 ‡ Moisture: not dec. 12 Date Analyzed: 01/02/90
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	11	U
74-83-9	-----Bromomethane	11	U
75-01-4	-----Vinyl Chloride	11	U
75-00-3	-----Chloroethane	11	U
75-09-2	-----Methylene Chloride	25	B
67-64-1	-----Acetone	61	
75-15-0	-----Carbon Disulfide	6	U
75-35-4	-----1,1-Dichloroethene	6	U
75-34-3	-----1,1-Dichloroethane	6	U
540-59-0	-----1,2-Dichloroethene (total)	6	U
67-66-3	-----Chloroform	6	U
107-06-2	-----1,2-Dichloroethane	6	U
78-93-3	-----2-Butanone	33	
71-55-6	-----1,1,1-Trichloroethane	6	U
56-23-5	-----Carbon Tetrachloride	6	U
108-05-4	-----Vinyl Acetate	11	U
75-27-4	-----Bromodichloromethane	6	U
78-87-5	-----1,2-Dichloropropane	6	U
10061-01-5	-----cis-1,3-Dichloropropene	6	U
79-01-6	-----Trichloroethene	6	U
124-48-1	-----Dibromochloromethane	6	U
79-00-5	-----1,1,2-Trichloroethane	6	U
71-43-2	-----Benzene	6	U
10061-02-6	-----Trans-1,3-Dichloropropene	6	U
75-25-2	-----Bromoform	6	U
108-10-1	-----4-Methyl-2-Pentanone	11	U
591-78-6	-----2-Hexanone	15	
127-18-4	-----Tetrachloroethene	6	U
79-34-5	-----1,1,2,2-Tetrachloroethane	6	U
108-88-3	-----Toluene	6	U
108-90-7	-----Chlorobenzene	6	U
100-41-4	-----Ethylbenzene	6	U
100-42-5	-----Styrene	6	U
1330-20-7	-----Total Xylenes	6	U

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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201BREMSD

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: 311272

Sample wt/vol: 4.0 (g/mL) G Lab File ID: CN011272B13

Level: (low/med) MED Date Received: 12/20/89

% Moisture: not dec. 12 Date Analyzed: 12/29/89

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	1400	U
74-83-9	-----Bromomethane	1400	U
75-01-4	-----Vinyl Chloride	1400	U
75-00-3	-----Chloroethane	1400	U
75-09-2	-----Methylene Chloride	530	BJ
67-64-1	-----Acetone	1400	U
75-15-0	-----Carbon Disulfide	710	U
75-35-4	-----1,1-Dichloroethene	710	U
75-34-3	-----1,1-Dichloroethane	710	U
540-59-0	-----1,2-Dichloroethene (total)	710	U
67-66-3	-----Chloroform	710	U
107-06-2	-----1,2-Dichloroethane	710	U
78-93-3	-----2-Butanone	1400	U
71-55-6	-----1,1,1-Trichloroethane	710	U
56-23-5	-----Carbon Tetrachloride	710	U
108-05-4	-----Vinyl Acetate	1400	U
75-27-4	-----Bromodichloromethane	710	U
78-87-5	-----1,2-Dichloropropane	710	U
10061-01-5	-----cis-1,3-Dichloropropene	710	U
79-01-6	-----Trichloroethene	710	U
124-48-1	-----Dibromochloromethane	710	U
79-00-5	-----1,1,2-Trichloroethane	710	U
71-43-2	-----Benzene	710	U
10061-02-6	-----Trans-1,3-Dichloropropene	710	U
75-25-2	-----Bromoform	710	U
108-10-1	-----4-Methyl-2-Pentanone	1400	U
591-78-6	-----2-Hexanone	3600	
127-18-4	-----Tetrachloroethene	710	U
79-34-5	-----1,1,2,2-Tetrachloroethane	710	U
108-88-3	-----Toluene	710	U
108-90-7	-----Chlorobenzene	710	U
100-41-4	-----Ethylbenzene	7700	
100-42-5	-----Styrene	2900	
1330-20-7	-----Total Xylenes	22000	

FORM I VOA

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1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G3D09679B08
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/28/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 20

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	44000	
111-44-4	bis(2-Chloroethyl) Ether	7300	U
95-57-8	2-Chlorophenol	7300	U
541-73-1	1,3-Dichlorobenzene	7300	U
106-46-7	1,4-Dichlorobenzene	7300	U
100-51-6	Benzyl Alcohol	7300	U
95-50-1	1,2-Dichlorobenzene	7300	U
95-48-7	2-Methylphenol	130000	E
39638-32-9	bis(2-Chloroisopropyl) Ether	7300	U
106-44-5	4-Methylphenol	180000	E
621-64-7	N-Nitroso-Di-n-Propylamine	7300	U
67-72-1	Hexachloroethane	7300	U
98-95-3	Nitrobenzene	7300	U
78-59-1	Isophorone	7300	U
88-75-5	2-Nitrophenol	7300	U
105-67-9	2,4-Dimethylphenol	240000	E
65-85-0	Benzoic Acid	35000	U
111-91-1	bis(2-Chloroethoxy) Methane	7300	U
120-83-2	2,4-Dichlorophenol	7300	U
120-82-1	1,2,4-Trichlorobenzene	7300	U
91-20-3	Naphthalene	54000	
106-47-8	4-Chloroaniline	7300	U
87-68-3	Hexachlorobutadiene	7300	U
59-50-7	4-Chloro-3-Methylphenol	7300	U
91-57-6	2-Methylnaphthalene	49000	
77-47-4	Hexachlorocyclopentadiene	7300	U
88-06-2	2,4,6-Trichlorophenol	7300	U
95-95-4	2,4,5-Trichlorophenol	35000	U
91-58-7	2-Chloronaphthalene	7300	U
88-74-4	2-Nitroaniline	35000	U
131-11-3	Dimethyl Phthalate	7300	U
208-96-8	Acenaphthylene	4200	J
606-20-2	2,6-Dinitrotoluene	7300	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G3D09679B08
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/28/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 20

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline	35000	U
83-32-9-----	Acenaphthene	3700	J
51-28-5-----	2,4-Dinitrophenol	35000	U
100-02-7-----	4-Nitrophenol	35000	U
132-64-9-----	Dibenzofuran	14000	
121-14-2-----	2,4-Dinitrotoluene	7300	U
84-66-2-----	Diethylphthalate	7300	U
7005-72-3-----	4-Chlorophenyl-phenylether	7300	U
86-73-7-----	Fluorene	7300	U
100-01-6-----	4-Nitroaniline	35000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	35000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	7300	U
101-55-3-----	4-Bromophenyl-phenylether	7300	U
118-74-1-----	Hexachlorobenzene	7300	U
87-86-5-----	Pentachlorophenol	35000	U
85-01-8-----	Phenanthrene	19000	
120-12-7-----	Anthracene	4700	J
84-74-2-----	Di-n-Butylphthalate	7300	U
206-44-0-----	Fluoranthene	3600	J
129-00-0-----	Pyrene	3200	J
85-68-7-----	Butylbenzylphthalate	7300	U
91-94-1-----	3,3'-Dichlorobenzidine	15000	U
56-55-3-----	Benzo(a)Anthracene	2300	J
218-01-9-----	Chrysene	1500	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate	7300	U
117-84-0-----	Di-n-Octyl Phthalate	7300	U
205-99-2-----	Benzo(b)Fluoranthene	780	JX
207-08-9-----	Benzo(k)Fluoranthene	780	JX
50-32-8-----	Benzo(a)Pyrene	7300	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	7300	U
53-70-3-----	Dibenzo(a,h)Anthracene	7300	U
191-24-2-----	Benzo(g,h,i)Perylene	7300	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

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1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G3D09679B08
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/28/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 20

Number TICs found: 20 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 28790-86-5	2-CYCLOPENTEN-1-ONE, 2,3,4-T	7.57	110000	J
2. 526-75-0	PHENOL, 2,3-DIMETHYL-	7.73	88000	J
3. 526-75-0	PHENOL, 2,3-DIMETHYL-	8.27	170000	J
4. 93-51-6	PHENOL, 2-METHOXY-4-METHYL-	8.44	88000	J
5. 28715-26-6	BENZOFURAN, 4,7-DIMETHYL-	8.80	67000	J
6. 698-71-5	PHENOL, 3-ETHYL-5-METHYL-	8.92	73000	J
7. 5873-86-9	BENZENECARBOETHIOIC ACID, O-M	9.19	56000	J
8. 24599-58-4	BENZENE, 1,4-DIMETHOXY-2-MET	9.32	150000	J
9.	UNKNOWN	9.64	36000	J
10. 91-57-6	NAPHTHALENE, 2-METHYL-	9.72	34000	J
11. 91-10-1	PHENOL, 2,6-DIMETHOXY-	9.90	95000	J
12.	UNKNOWN	9.99	45000	J
13.	UNKNOWN	10.07	59000	J
14. 17453-94-0	UNDECANE, 5-ETHYL-	10.24	47000	J
15. 121-34-6	BENZOIC ACID, 4-HYDROXY-3-ME	10.69	190000	J
16.	UNKNOWN	11.29	130000	J
17. 54105-67-8	HEPTADECANE, 2,6-DIMETHYL-	11.74	48000	J
18.	UNKNOWN	11.89	95000	J
19. 629-78-7	HEPTADECANE	12.44	56000	J
20. 544-76-3	HEXADECANE	13.10	37000	J

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201ADL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G2D09679C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 50

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	23000	D
111-44-4	bis(2-Chloroethyl) Ether	18000	U
95-57-8	2-Chlorophenol	18000	U
541-73-1	1,3-Dichlorobenzene	18000	U
106-46-7	1,4-Dichlorobenzene	18000	U
100-51-6	Benzyl Alcohol	18000	U
95-50-1	1,2-Dichlorobenzene	18000	U
95-48-7	2-Methylphenol	120000	D
39638-32-9	bis(2-Chloroisopropyl) Ether	18000	U
106-44-5	4-Methylphenol	170000	D
621-64-7	N-Nitroso-Di-n-Propylamine	18000	U
67-72-1	Hexachloroethane	18000	U
98-95-3	Nitrobenzene	18000	U
78-59-1	Isophorone	18000	U
88-75-5	2-Nitrophenol	18000	U
105-67-9	2,4-Dimethylphenol	170000	D
65-85-0	Benzoic Acid	88000	U
111-91-1	bis(2-Chloroethoxy) Methane	18000	U
120-83-2	2,4-Dichlorophenol	18000	U
120-82-1	1,2,4-Trichlorobenzene	18000	U
91-20-3	Naphthalene	42000	D
106-47-8	4-Chloroaniline	18000	U
87-68-3	Hexachlorobutadiene	18000	U
59-50-7	4-Chloro-3-Methylphenol	18000	U
91-57-6	2-Methylnaphthalene	27000	D
77-47-4	Hexachlorocyclopentadiene	18000	U
88-06-2	2,4,6-Trichlorophenol	18000	U
95-95-4	2,4,5-Trichlorophenol	88000	U
91-58-7	2-Chloronaphthalene	18000	U
88-74-4	2-Nitroaniline	88000	U
131-11-3	Dimethyl Phthalate	18000	U
208-96-8	Acenaphthylene	3600	DJ
606-20-2	2,6-Dinitrotoluene	18000	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201ADL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G2D09679C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 50

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

99-09-2-----3-Nitroaniline_____	88000	U
83-32-9-----Acenaphthene_____	2600	DJ
51-28-5-----2,4-Dinitrophenol_____	88000	U
100-02-7-----4-Nitrophenol_____	88000	U
132-64-9-----Dibenzofuran_____	7900	DJ
121-14-2-----2,4-Dinitrotoluene_____	18000	U
84-66-2-----Diethylphthalate_____	18000	U
7005-72-3-----4-Chlorophenyl-phenylether_____	18000	U
86-73-7-----Fluorene_____	18000	U
100-01-6-----4-Nitroaniline_____	88000	U
534-52-1-----4,6-Dinitro-2-Methylphenol_____	88000	U
86-30-6-----N-Nitrosodiphenylamine (1)_____	18000	U
101-55-3-----4-Bromophenyl-phenylether_____	18000	U
118-74-1-----Hexachlorobenzene_____	18000	U
87-86-5-----Pentachlorophenol_____	88000	U
85-01-8-----Phenanthrene_____	7700	DJ
120-12-7-----Anthracene_____	1900	DJ
84-74-2-----Di-n-Butylphthalate_____	18000	U
206-44-0-----Fluoranthene_____	2100	DJ
129-00-0-----Pyrene_____	2900	DJ
85-68-7-----Butylbenzylphthalate_____	18000	U
91-94-1-----3,3'-Dichlorobenzidine_____	36000	U
56-55-3-----Benzo(a)Anthracene_____	18000	U
218-01-9-----Chrysene_____	18000	U
117-81-7-----bis(2-Ethylhexyl) Phthalate_____	18000	U
117-84-0-----Di-n-Octyl Phthalate_____	18000	U
205-99-2-----Benzo(b) Fluoranthene_____	18000	U
207-08-9-----Benzo(k) Fluoranthene_____	18000	U
50-32-8-----Benzo(a) Pyrene_____	18000	U
193-39-5-----Indeno(1,2,3-cd) Pyrene_____	18000	U
53-70-3-----Dibenzo(a,h) Anthracene_____	18000	U
191-24-2-----Benzo(g,h,i) Perylene_____	18000	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201ADL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G2D09679C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 50

Number TICs found: 20 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	7.07	110000	J
2.	DIMETHYLPHENOL	7.23	100000	J
3.	DIMETHYLPHENOL	7.73	140000	J
4.	DIMETHYLPHENOL	7.83	72000	J
5.	UNKNOWN SUBSTITUTED HYDROCAR	8.23	55000	J
6.	UNKNOWN SUBSTITUTED HYDROCAR	8.34	61000	J
7.	UNKNOWN	8.59	84000	J
8.	UNKNOWN SUBSTITUTED HYDROCAR	8.72	290000	J
9.	UNKNOWN	9.00	46000	J
10.	UNKNOWN	9.07	48000	J
11.	DIMETHOXY PHENOL	9.25	150000	J
12.	UNKNOWN	9.34	57000	J
13.	UNKNOWN	9.40	84000	J
14.	UNKNOWN	9.57	42000	J
15.	UNKNOWN	9.77	48000	J
16.	UNKNOWN	9.99	350000	J
17.	UNKNOWN	10.54	280000	J
18.	UNKNOWN	11.10	180000	J
19.	UNKNOWN	11.57	44000	J
20.	UNKNOWN HYDROCARBON	12.22	26000	J

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GD009686A21
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 12 dec. _____ Date Extracted: 12/29/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/02/90
 GPC Cleanup: (Y/N) N pH: 5.8 Dilution Factor: 5.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2-----	Phenol_____	120000	
111-44-4-----	bis(2-Chloroethyl)Ether_____	110000	U
95-57-8-----	2-Chlorophenol_____	110000	U
541-73-1-----	1,3-Dichlorobenzene_____	110000	U
106-46-7-----	1,4-Dichlorobenzene_____	110000	U
100-51-6-----	Benzyl Alcohol_____	110000	U
95-50-1-----	1,2-Dichlorobenzene_____	110000	U
95-48-7-----	2-Methylphenol_____	400000	
39638-32-9-----	bis(2-Chloroisopropyl)Ether_____	110000	U
106-44-5-----	4-Methylphenol_____	690000	
621-64-7-----	N-Nitroso-Di-n-Propylamine_____	110000	U
67-72-1-----	Hexachloroethane_____	110000	U
98-95-3-----	Nitrobenzene_____	110000	U
78-59-1-----	Isophorone_____	110000	U
88-75-5-----	2-Nitrophenol_____	110000	U
105-67-9-----	2,4-Dimethylphenol_____	610000	
65-85-0-----	Benzoic Acid_____	550000	U
111-91-1-----	bis(2-Chloroethoxy)Methane_____	110000	U
120-83-2-----	2,4-Dichlorophenol_____	110000	U
120-82-1-----	1,2,4-Trichlorobenzene_____	110000	U
91-20-3-----	Naphthalene_____	160000	
106-47-8-----	4-Chloroaniline_____	110000	U
87-68-3-----	Hexachlorobutadiene_____	110000	U
59-50-7-----	4-Chloro-3-Methylphenol_____	110000	U
91-57-6-----	2-Methylnaphthalene_____	120000	
77-47-4-----	Hexachlorocyclopentadiene_____	110000	U
88-06-2-----	2,4,6-Trichlorophenol_____	110000	U
95-95-4-----	2,4,5-Trichlorophenol_____	550000	U
91-58-7-----	2-Chloronaphthalene_____	110000	U
88-74-4-----	2-Nitroaniline_____	550000	U
131-11-3-----	Dimethyl Phthalate_____	110000	U
208-96-8-----	Acenaphthylene_____	11000	J
606-20-2-----	2,6-Dinitrotoluene_____	110000	U

FORM I SV-1

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: 309686

Sample wt/vol: 1.0 (g/mL) G Lab File ID: GD009686A21

Level: (low/med) MED Date Received: 12/20/89

% Moisture: not dec. 12 dec. _____ Date Extracted: 12/29/89

Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/02/90

GPC Cleanup: (Y/N) N pH: 5.8 Dilution Factor: 5.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline	550000	U
83-32-9-----	Acenaphthene	110000	U
51-28-5-----	2,4-Dinitrophenol	550000	U
100-02-7-----	4-Nitrophenol	550000	U
132-64-9-----	Dibenzofuran	43000	J
121-14-2-----	2,4-Dinitrotoluene	110000	U
84-66-2-----	Diethylphthalate	110000	U
7005-72-3-----	4-Chlorophenyl-phenylether	110000	U
86-73-7-----	Fluorene	35000	J
100-01-6-----	4-Nitroaniline	550000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	550000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	110000	U
101-55-3-----	4-Bromophenyl-phenylether	110000	U
118-74-1-----	Hexachlorobenzene	110000	U
87-86-5-----	Pentachlorophenol	550000	U
85-01-8-----	Phenanthrene	46000	J
120-12-7-----	Anthracene	11000	J
84-74-2-----	Di-n-Butylphthalate	110000	U
206-44-0-----	Fluoranthene	110000	U
129-00-0-----	Pyrene	14000	J
85-68-7-----	Butylbenzylphthalate	110000	U
91-94-1-----	3,3'-Dichlorobenzidine	230000	U
56-55-3-----	Benzo(a)Anthracene	110000	U
218-01-9-----	Chrysene	110000	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	110000	U
117-84-0-----	Di-n-Octyl Phthalate	110000	U
205-99-2-----	Benzo(b) Fluoranthene	110000	U
207-08-9-----	Benzo(k) Fluoranthene	110000	U
50-32-8-----	Benzo(a) Pyrene	110000	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene	110000	U
53-70-3-----	Dibenzo(a,h) Anthracene	110000	U
191-24-2-----	Benzo(g,h,i) Perylene	110000	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GD009686A21
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 12 dec. _____ Date Extracted: 12/29/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/02/90
 GPC Cleanup: (Y/N) N pH: 5.8 Dilution Factor: 5.0

Number TICs found: 20 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	8.22	500000	J
2.	104-55-2 2-PROPENAL, 3-PHENYL-	8.45	34000	J
3.	UNKNOWN DIMETHYL PHENOL	8.87	140000	J
4.	93-51-6 PHENOL, 2-METHOXY-4-METHYL-	9.05	45000	J
5.	UNKNOWN DIMETHYL PHENOL	9.10	68000	J
6.	UNKNOWN	9.44	68000	J
7.	UNKNOWN	9.50	80000	J
8.	UNKNOWN TRIMETHYL PHENOL	9.80	100000	J
9.	UNKNOWN SUBST. HYDROCARBON	9.92	320000	J
10.	UNKNOWN	10.34	57000	J
11.	UNKNOWN SUBST. HYDROCARBON	10.47	170000	J
12.	UNKNOWN	10.55	68000	J
13.	UNKNOWN	10.64	110000	J
14.	UNKNOWN	10.82	68000	J
15.	UNKNOWN	11.22	310000	J
16.	UNKNOWN	11.79	330000	J
17.	UNKNOWN	12.22	45000	J
18.	UNKNOWN	12.37	190000	J
19.	UNKNOWN HYDROCARBON	12.89	80000	J
20.	UNKNOWN HYDROCARBON	13.50	91000	J

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309688
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009688C20
 Level: (low/med) LOW Date Received: 12/20/89
 ‡ Moisture: not dec. 12 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 8.1 Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	370	U
111-44-4	bis(2-Chloroethyl) Ether	370	U
95-57-8	2-Chlorophenol	370	U
541-73-1	1,3-Dichlorobenzene	370	U
106-46-7	1,4-Dichlorobenzene	370	U
100-51-6	Benzyl Alcohol	370	U
95-50-1	1,2-Dichlorobenzene	370	U
95-48-7	2-Methylphenol	52	J
39638-32-9	bis(2-Chloroisopropyl) Ether	370	U
106-44-5	4-Methylphenol	6000	
621-64-7	N-Nitroso-Di-n-Propylamine	370	U
67-72-1	Hexachloroethane	370	U
98-95-3	Nitrobenzene	370	U
78-59-1	Isophorone	370	U
88-75-5	2-Nitrophenol	370	U
105-67-9	2,4-Dimethylphenol	1600	
65-85-0	Benzoic Acid	1800	U
111-91-1	bis(2-Chloroethoxy) Methane	370	U
120-83-2	2,4-Dichlorophenol	370	U
120-82-1	1,2,4-Trichlorobenzene	370	U
91-20-3	Naphthalene	67	J
106-47-8	4-Chloroaniline	370	U
87-68-3	Hexachlorobutadiene	370	U
59-50-7	4-Chloro-3-Methylphenol	370	U
91-57-6	2-Methylnaphthalene	80	J
77-47-4	Hexachlorocyclopentadiene	370	U
88-06-2	2,4,6-Trichlorophenol	370	U
95-95-4	2,4,5-Trichlorophenol	1800	U
91-58-7	2-Chloronaphthalene	370	U
88-74-4	2-Nitroaniline	1800	U
131-11-3	Dimethyl Phthalate	370	U
208-96-8	Acenaphthylene	370	U
606-20-2	2,6-Dinitrotoluene	370	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: 309688

Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009688C20

Level: (low/med) LOW Date Received: 12/20/89

% Moisture: not dec. 12 dec. _____ Date Extracted: 12/21/89

Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89

GPC Cleanup: (Y/N) N pH: 8.1 Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline	1800	U
83-32-9-----	Acenaphthene	370	U
51-28-5-----	2,4-Dinitrophenol	1800	U
100-02-7-----	4-Nitrophenol	1800	U
132-64-9-----	Dibenzofuran	370	U
121-14-2-----	2,4-Dinitrotoluene	370	U
84-66-2-----	Diethylphthalate	370	U
7005-72-3-----	4-Chlorophenyl-phenylether	370	U
86-73-7-----	Fluorene	74	J
100-01-6-----	4-Nitroaniline	1800	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1800	U
86-30-6-----	N-Nitrosodiphenylamine (1)	370	U
101-55-3-----	4-Bromophenyl-phenylether	370	U
118-74-1-----	Hexachlorobenzene	370	U
87-86-5-----	Pentachlorophenol	1800	U
85-01-8-----	Phenanthrene	39	J
120-12-7-----	Anthracene	370	U
84-74-2-----	Di-n-Butylphthalate	370	U
206-44-0-----	Fluoranthene	370	U
129-00-0-----	Pyrene	370	U
85-68-7-----	Butylbenzylphthalate	370	U
91-94-1-----	3,3'-Dichlorobenzidine	750	U
56-55-3-----	Benzo(a)Anthracene	370	U
218-01-9-----	Chrysene	370	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	63	J
117-84-0-----	Di-n-Octyl Phthalate	370	U
205-99-2-----	Benzo(b)Fluoranthene	370	U
207-08-9-----	Benzo(k)Fluoranthene	370	U
50-32-8-----	Benzo(a)Pyrene	370	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	370	U
53-70-3-----	Dibenzo(a,h)Anthracene	370	U
191-24-2-----	Benzo(g,h,i)Perylene	370	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309688
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009688C20
 Level: (low/med) LOW Date Received: 12/20/89
 ‡ Moisture: not dec. 12 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 8.1 Dilution Factor: 1.00

Number TICs found: 11 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	6.60	190	J
2.	DIMETHYLPHENOL ISOMER	7.23	300	J
3.	DIMETHYLPHENOL ISOMER	7.47	150	J
4.	DIMETHYLPHENOL ISOMER	7.75	2700	J
5.	DIMETHYLPHENOL ISOMER	7.83	190	J
6. 621-27-2	PHENOL, 3-PROPYL-	8.50	1200	J
7. 2416-94-6	PHENOL, 2,3,6-TRIMETHYL-	8.64	340	J
8.	UNKNOWN	9.20	190	J
9.	UNKNOWN	9.35	530	J
10.	UNKNOWN	10.04	340	J
11.	UNKNOWN	11.10	230	J

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009689C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 14 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 7.7 Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
108-95-2	Phenol	1300	
111-44-4	bis(2-Chloroethyl) Ether	380	U
95-57-8	2-Chlorophenol	380	U
541-73-1	1,3-Dichlorobenzene	380	U
106-46-7	1,4-Dichlorobenzene	380	U
100-51-6	Benzyl Alcohol	380	U
95-50-1	1,2-Dichlorobenzene	380	U
95-48-7	2-Methylphenol	2400	
39638-32-9	bis(2-Chloroisopropyl) Ether	380	U
106-44-5	4-Methylphenol	13000	E
621-64-7	N-Nitroso-Di-n-Propylamine	380	U
67-72-1	Hexachloroethane	380	U
98-95-3	Nitrobenzene	380	U
78-59-1	Isophorone	380	U
88-75-5	2-Nitrophenol	380	U
105-67-9	2,4-Dimethylphenol	7300	E
65-85-0	Benzoic Acid	1900	U
111-91-1	bis(2-Chloroethoxy) Methane	380	U
120-83-2	2,4-Dichlorophenol	380	U
120-82-1	1,2,4-Trichlorobenzene	380	U
91-20-3	Naphthalene	1200	
106-47-8	4-Chloroaniline	380	U
87-68-3	Hexachlorobutadiene	380	U
59-50-7	4-Chloro-3-Methylphenol	380	U
91-57-6	2-Methylnaphthalene	1800	
77-47-4	Hexachlorocyclopentadiene	380	U
88-06-2	2,4,6-Trichlorophenol	380	U
95-95-4	2,4,5-Trichlorophenol	1900	U
91-58-7	2-Chloronaphthalene	380	U
88-74-4	2-Nitroaniline	1900	U
131-11-3	Dimethyl Phthalate	380	U
208-96-8	Acenaphthylene	140	J
606-20-2	2,6-Dinitrotoluene	380	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009689C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 14 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 7.7 Dilution Factor: 1.00

CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline_____	1900	U
83-32-9-----	Acenaphthene_____	180	J
51-28-5-----	2,4-Dinitrophenol_____	1900	U
100-02-7-----	4-Nitrophenol_____	1900	U
132-64-9-----	Dibenzofuran_____	600	
121-14-2-----	2,4-Dinitrotoluene_____	380	U
84-66-2-----	Diethylphthalate_____	380	U
7005-72-3-----	4-Chlorophenyl-phenylether_____	380	U
86-73-7-----	Fluorene_____	570	
100-01-6-----	4-Nitroaniline_____	1900	U
534-52-1-----	4,6-Dinitro-2-Methylphenol_____	1900	U
86-30-6-----	N-Nitrosodiphenylamine (1)_____	380	U
101-55-3-----	4-Bromophenyl-phenylether_____	380	U
118-74-1-----	Hexachlorobenzene_____	380	U
87-86-5-----	Pentachlorophenol_____	1900	U
85-01-8-----	Phenanthrene_____	860	
120-12-7-----	Anthracene_____	180	J
84-74-2-----	Di-n-Butylphthalate_____	82	J
206-44-0-----	Fluoranthene_____	230	J
129-00-0-----	Pyrene_____	270	J
85-68-7-----	Butylbenzylphthalate_____	380	U
91-94-1-----	3,3'-Dichlorobenzidine_____	770	U
56-55-3-----	Benzo(a)Anthracene_____	63	J
218-01-9-----	Chrysene_____	74	J
117-81-7-----	bis(2-Ethylhexyl)Phthalate_____	88	J
117-84-0-----	Di-n-Octyl Phthalate_____	380	U
205-99-2-----	Benzo(b)Fluoranthene_____	380	U
207-08-9-----	Benzo(k)Fluoranthene_____	380	U
50-32-8-----	Benzo(a)Pyrene_____	380	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene_____	380	U
53-70-3-----	Dibenzo(a,h)Anthracene_____	380	U
191-24-2-----	Benzo(g,h,i)Perylene_____	380	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009689C20
 Level: (low/med) LOW Date Received: 12/20/89
 ‡ Moisture: not dec. 14 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 7.7 Dilution Factor: 1.00

Number TICs found: 20

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	DIMETHYLPHENOL + ISOMER	7.75	2800	J
2. 621-27-2	PHENOL, 3-PROPYL-	8.50	1500	J
3.	UNKNOWN	9.37	1300	J
4.	UNKNOWN	9.57	1300	J
5.	DIMETHYLNAPHTHALENE ISOMER	9.90	1400	J
6.	UNKNOWN	10.05	1500	J
7. 54105-67-8	HEPTADECANE, 2,6-DIMETHYL-	10.95	1900	J
8.	UNKNOWN	11.10	1000	J
9. 629-78-7	HEPTADECANE	11.59	1500	J
10. 54105-67-8	HEPTADECANE, 2,6-DIMETHYL-	12.20	1100	J
11.	UNKNOWN	12.55	1100	J
12. 544-76-3	HEXADECANE	12.77	1800	J
13. 57-10-3	HEXADECANOIC ACID	13.12	1400	J
14.	HEPTADECANE ISOMER	13.32	1600	J
15.	PENTACOSANE ISOMER	13.84	2300	J
16.	PENTACOSANE ISOMER	14.80	2400	J
17. 7206-19-1	3-OCTADECENE, (E)-	15.24	5000	J
18.	UNKNOWN	15.70	1400	J
19.	UNKNOWN	16.12	4600	J
20.	UNKNOWN	16.60	2100	J

FORM I SV-TIC

1/87 Rev.

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202BDL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GD009689A20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 14 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 7.7 Dilution Factor: 5.0

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
108-95-2	Phenol	920	DJ
111-44-4	bis(2-Chloroethyl) Ether	1900	U
95-57-8	2-Chlorophenol	1900	U
541-73-1	1,3-Dichlorobenzene	1900	U
106-46-7	1,4-Dichlorobenzene	1900	U
100-51-6	Benzyl Alcohol	1900	U
95-50-1	1,2-Dichlorobenzene	1900	U
95-48-7	2-Methylphenol	2200	D
39638-32-9	bis(2-Chloroisopropyl) Ether	1900	U
106-44-5	4-Methylphenol	12000	D
621-64-7	N-Nitroso-Di-n-Propylamine	1900	U
67-72-1	Hexachloroethane	1900	U
98-95-3	Nitrobenzene	1900	U
78-59-1	Isophorone	1900	U
88-75-5	2-Nitrophenol	1900	U
105-67-9	2,4-Dimethylphenol	7300	D
65-85-0	Benzoic Acid	9300	U
111-91-1	bis(2-Chloroethoxy) Methane	1900	U
120-83-2	2,4-Dichlorophenol	1900	U
120-82-1	1,2,4-Trichlorobenzene	1900	U
91-20-3	Naphthalene	1300	DJ
106-47-8	4-Chloroaniline	1900	U
87-68-3	Hexachlorobutadiene	1900	U
59-50-7	4-Chloro-3-Methylphenol	1900	U
91-57-6	2-Methylnaphthalene	1800	DJ
77-47-4	Hexachlorocyclopentadiene	1900	U
88-06-2	2,4,6-Trichlorophenol	1900	U
95-95-4	2,4,5-Trichlorophenol	9300	U
91-58-7	2-Chloronaphthalene	1900	U
88-74-4	2-Nitroaniline	9300	U
131-11-3	Dimethyl Phthalate	1900	U
208-96-8	Acenaphthylene	1900	U
606-20-2	2,6-Dinitrotoluene	1900	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202BDL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GD009689A20
 Level: (low/med) LOW Date Received: 12/20/89
 ‡ Moisture: not dec. 14 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 7.7 Dilution Factor: 5.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2	3-Nitroaniline	9300	U
83-32-9	Acenaphthene	220	DJ
51-28-5	2,4-Dinitrophenol	9300	U
100-02-7	4-Nitrophenol	9300	U
132-64-9	Dibenzofuran	660	DJ
121-14-2	2,4-Dinitrotoluene	1900	U
84-66-2	Diethylphthalate	1900	U
7005-72-3	4-Chlorophenyl-phenylether	1900	U
86-73-7	Fluorene	580	DJ
100-01-6	4-Nitroaniline	9300	U
534-52-1	4,6-Dinitro-2-Methylphenol	9300	U
86-30-6	N-Nitrosodiphenylamine (1)	1900	U
101-55-3	4-Bromophenyl-phenylether	1900	U
118-74-1	Hexachlorobenzene	1900	U
87-86-5	Pentachlorophenol	9300	U
85-01-8	Phenanthrene	900	DJ
120-12-7	Anthracene	200	DJ
84-74-2	Di-n-Butylphthalate	200	DJ
206-44-0	Fluoranthene	260	DJ
129-00-0	Pyrene	260	DJ
85-68-7	Butylbenzylphthalate	1900	U
91-94-1	3,3'-Dichlorobenzidine	3800	U
56-55-3	Benzo(a)Anthracene	1900	U
218-01-9	Chrysene	1900	U
117-81-7	bis(2-Ethylhexyl) Phthalate	1900	U
117-84-0	Di-n-Octyl Phthalate	1900	U
205-99-2	Benzo(b)Fluoranthene	1900	U
207-08-9	Benzo(k)Fluoranthene	1900	U
50-32-8	Benzo(a)Pyrene	1900	U
193-39-5	Indeno(1,2,3-cd)Pyrene	1900	U
53-70-3	Dibenzo(a,h)Anthracene	1900	U
191-24-2	Benzo(g,h,i)Perylene	1900	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202BDL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GD009689A20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 14 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 7.7 Dilution Factor: 5.0

Number TICs found: 20 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	DIMETHYLPHENOL ISOMER	7.70	6400	J
2. 621-27-2	PHENOL, 3-PROPYL-	8.47	3500	J
3. 2416-94-6	PHENOL, 2,3,6-TRIMETHYL-	8.59	1700	J
4.	UNKNOWN	8.97	1600	J
5. 4453-90-1	1,4-METHANONAPHTHALENE, 1,4-	9.05	1600	J
6.	UNKNOWN	9.54	2100	J
7. 569-41-5	NAPHTHALENE, 1,8-DIMETHYL-	9.85	2100	J
8.	UNKNOWN	10.00	2100	J
9. 54105-67-8	HEPTADECANE, 2,6-DIMETHYL-	10.92	2500	J
10. 629-78-7	HEPTADECANE	11.55	3100	J
11. 54105-67-8	HEPTADECANE, 2,6-DIMETHYL-	12.15	2300	J
12.	UNKNOWN	12.50	1900	J
13. 544-76-3	HEXADECANE	12.70	3100	J
14.	PENTACOSANE ISOMER	13.25	2700	J
15.	PENTACOSANE ISOMER	13.77	3900	J
16. 629-97-0	DOCOSANE	14.27	8700	J
17.	PENTACOSANE ISOMER	14.75	3100	J
18.	UNKNOWN	15.19	6000	J
19.	UNKNOWN	16.04	6200	J
20.	UNKNOWN	16.50	2900	J

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1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202C

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309690
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009690A20
 Level: (low/med) LOW Date Received: 12/20/89
 ‡ Moisture: not dec. 13 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 8.2 Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	380	U
111-44-4	bis(2-Chloroethyl) Ether	380	U
95-57-8	2-Chlorophenol	380	U
541-73-1	1,3-Dichlorobenzene	380	U
106-46-7	1,4-Dichlorobenzene	380	U
100-51-6	Benzyl Alcohol	380	U
95-50-1	1,2-Dichlorobenzene	380	U
95-48-7	2-Methylphenol	380	U
39638-32-9	bis(2-Chloroisopropyl) Ether	380	U
106-44-5	4-Methylphenol	510	
621-64-7	N-Nitroso-Di-n-Propylamine	380	U
67-72-1	Hexachloroethane	380	U
98-95-3	Nitrobenzene	380	U
78-59-1	Isophorone	380	U
88-75-5	2-Nitrophenol	380	U
105-67-9	2,4-Dimethylphenol	1400	
65-85-0	Benzoic Acid	1800	U
111-91-1	bis(2-Chloroethoxy) Methane	380	U
120-83-2	2,4-Dichlorophenol	380	U
120-82-1	1,2,4-Trichlorobenzene	380	U
91-20-3	Naphthalene	130	J
106-47-8	4-Chloroaniline	380	U
87-68-3	Hexachlorobutadiene	380	U
59-50-7	4-Chloro-3-Methylphenol	380	U
91-57-6	2-Methylnaphthalene	140	J
77-47-4	Hexachlorocyclopentadiene	380	U
88-06-2	2,4,6-Trichlorophenol	380	U
95-95-4	2,4,5-Trichlorophenol	1800	U
91-58-7	2-Chloronaphthalene	380	U
88-74-4	2-Nitroaniline	1800	U
131-11-3	Dimethyl Phthalate	380	U
208-96-8	Acenaphthylene	380	U
606-20-2	2,6-Dinitrotoluene	380	U

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1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202C

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309690
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009690A20
 Level: (low/meg) LOW Date Received: 12/20/89
 % Moisture: not dec. 13 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 8.2 Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2	3-Nitroaniline	1800	U
83-32-9	Acenaphthene	380	U
51-28-5	2,4-Dinitrophenol	1800	U
100-02-7	4-Nitrophenol	1800	U
132-64-9	Dibenzofuran	39	J
121-14-2	2,4-Dinitrotoluene	380	U
84-66-2	Diethylphthalate	380	U
7005-72-3	4-Chlorophenyl-phenylether	380	U
86-73-7	Fluorene	59	J
100-01-6	4-Nitroaniline	1800	U
534-52-1	4,6-Dinitro-2-Methylphenol	1800	U
86-30-6	N-Nitrosodiphenylamine (1)	380	U
101-55-3	4-Bromophenyl-phenylether	380	U
118-74-1	Hexachlorobenzene	380	U
87-86-5	Pentachlorophenol	1800	U
85-01-8	Phenanthrene	380	U
120-12-7	Anthracene	380	U
84-74-2	Di-n-Butylphthalate	280	J
206-44-0	Fluoranthene	380	U
129-00-0	Pyrene	380	U
85-68-7	Butylbenzylphthalate	380	U
91-94-1	3,3'-Dichlorobenzidine	760	U
56-55-3	Benzo(a)Anthracene	380	U
218-01-9	Chrysene	380	U
117-81-7	bis(2-Ethylhexyl)Phthalate	77	J
117-84-0	Di-n-Octyl Phthalate	380	U
205-99-2	Benzo(b)Fluoranthene	380	U
207-08-9	Benzo(k)Fluoranthene	380	U
50-32-8	Benzo(a)Pyrene	380	U
193-39-5	Indeno(1,2,3-cd)Pyrene	380	U
53-70-3	Dibenzo(a,h)Anthracene	380	U
191-24-2	Benzo(g,h,i)Perylene	380	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202C

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309690
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH009690A20
 Level: (low/med) LOW Date Received: 12/20/89
 ‡ Moisture: not dec. 13 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 8.2 Dilution Factor: 1.00

Number TICs found: 13 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	ALDOL	5.62	380	AJ
2.	DIMETHYLPHENOL + UNKNOWN	7.25	310	J
3.	DIMETHYLPHENOL ISOMER	7.75	1000	J
4.	DIMETHYLPHENOL ISOMER	7.85	150	J
5.	TRIMETHYLPHENOL ISOMER	8.10	110	J
6.	ETHYLMETHYLPHENOL ISOMER	8.25	310	J
7.	TRIMETHYLPHENOL ISOMER	8.35	310	J
8.	ETHYLMETHYLPHENOL ISOMER	8.52	840	J
9.	TRIMETHYLPHENOL ISOMER	8.64	460	J
10.	UNKNOWN	9.00	230	J
11.	UNKNOWN	9.09	190	J
12.	UNKNOWN	9.20	270	J
13.	UNKNOWN	9.30	540	J

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TAR

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GRD09687A08
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: 8.5 Dilution Factor: 5.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	330000	
111-44-4	bis(2-Chloroethyl) Ether	280000	U
95-57-8	2-Chlorophenol	280000	U
541-73-1	1,3-Dichlorobenzene	280000	U
106-46-7	1,4-Dichlorobenzene	280000	U
100-51-6	Benzyl Alcohol	280000	U
95-50-1	1,2-Dichlorobenzene	280000	U
95-48-7	2-Methylphenol	1100000	
39638-32-9	bis(2-Chloroisopropyl) Ether	280000	U
106-44-5	4-Methylphenol	1400000	
621-64-7	N-Nitroso-Di-n-Propylamine	280000	U
67-72-1	Hexachloroethane	280000	U
98-95-3	Nitrobenzene	280000	U
78-59-1	Isophorone	280000	U
88-75-5	2-Nitrophenol	280000	U
105-67-9	2,4-Dimethylphenol	2000000	
65-85-0	Benzoic Acid	1300000	U
111-91-1	bis(2-Chloroethoxy) Methane	280000	U
120-83-2	2,4-Dichlorophenol	280000	U
120-82-1	1,2,4-Trichlorobenzene	280000	U
91-20-3	Naphthalene	340000	
106-47-8	4-Chloroaniline	280000	U
87-68-3	Hexachlorobutadiene	280000	U
59-50-7	4-Chloro-3-Methylphenol	280000	U
91-57-6	2-Methylnaphthalene	560000	
77-47-4	Hexachlorocyclopentadiene	280000	U
88-06-2	2,4,6-Trichlorophenol	280000	U
95-95-4	2,4,5-Trichlorophenol	1300000	U
91-58-7	2-Chloronaphthalene	280000	U
88-74-4	2-Nitroaniline	1300000	U
131-11-3	Dimethyl Phthalate	280000	U
208-96-8	Acenaphthylene	280000	U
606-20-2	2,6-Dinitrotoluene	280000	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TAR

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GRD09687A08
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: 8.5 Dilution Factor: 5.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline_____	1300000	U
83-32-9-----	Acenaphthene_____	280000	U
51-28-5-----	2,4-Dinitrophenol_____	1300000	U
100-02-7-----	4-Nitrophenol_____	1300000	U
132-64-9-----	Dibenzofuran_____	51000	J
121-14-2-----	2,4-Dinitrotoluene_____	280000	U
84-66-2-----	Diethylphthalate_____	280000	U
7005-72-3-----	4-Chlorophenyl-phenylether_____	280000	U
86-73-7-----	Fluorene_____	100000	J
100-01-6-----	4-Nitroaniline_____	1300000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol_____	1300000	U
86-30-6-----	N-Nitrosodiphenylamine (1)_____	280000	U
101-55-3-----	4-Bromophenyl-phenylether_____	280000	U
118-74-1-----	Hexachlorobenzene_____	280000	U
87-86-5-----	Pentachlorophenol_____	1300000	U
85-01-8-----	Phenanthrene_____	280000	U
120-12-7-----	Anthracene_____	280000	U
84-74-2-----	Di-n-Butylphthalate_____	280000	U
206-44-0-----	Fluoranthene_____	280000	U
129-00-0-----	Pyrene_____	280000	U
85-68-7-----	Butylbenzylphthalate_____	280000	U
91-94-1-----	3,3'-Dichlorobenzidine_____	550000	U
56-55-3-----	Benzo(a)Anthracene_____	280000	U
218-01-9-----	Chrysene_____	280000	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate_____	280000	U
117-84-0-----	Di-n-Octyl Phthalate_____	280000	U
205-99-2-----	Benzo(b)Fluoranthene_____	280000	U
207-08-9-----	Benzo(k)Fluoranthene_____	280000	U
50-32-8-----	Benzo(a)Pyrene_____	280000	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene_____	280000	U
53-70-3-----	Dibenzo(a,h)Anthracene_____	280000	U
191-24-2-----	Benzo(g,h,i)Perylene_____	280000	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202TAR

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GRD09687A08
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: 8.5 Dilution Factor: 5.0

Number TICs found: 20 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 620-14-4	BENZENE, 1-ETHYL-3-METHYL-	7.13	1100000	J
2.	UNKNOWN	8.35	1800000	J
3. 104-55-2	2-PROPENAL, 3-PHENYL-	8.60	1200000	J
4.	UNKNOWN	8.99	1200000	J
5.	UNKNOWN	9.20	1900000	J
6. 28715-26-6	BENZOFURAN, 4,7-DIMETHYL-	9.52	1400000	J
7.	UNKNOWN	9.84	1300000	J
8. 24599-58-4	BENZENE, 1,4-DIMETHOXY-2-MET	9.99	3300000	J
9.	DIMETHOXYPHENOL + UNKNOWN	10.52	1200000	J
10.	UNKNOWN	10.67	1200000	J
11.	DODECANE + UNKNOWN	10.75	1100000	J
12.	UNKNOWN	11.22	1500000	J
13. 544-76-3	HEXADECANE	11.45	1700000	J
14.	UNKNOWN	11.77	3300000	J
15.	UNKNOWN	12.00	1000000	J
16. 544-76-3	HEXADECANE	12.12	1600000	J
17. 629-78-7	HEPTADECANE	12.77	2100000	J
18. 54105-67-8	HEPTADECANE, 2,6-DIMETHYL-	13.37	1200000	J
19. 544-76-3	HEXADECANE	13.94	1100000	J
20.	UNKNOWN	16.60	3300000	J

FORM I SV-TIC

1/87 Rev.

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201ADLMS

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309683
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GD009683C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 50

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2-----	Phenol	18000	U
111-44-4-----	bis(2-Chloroethyl) Ether	18000	U
95-57-8-----	2-Chlorophenol	18000	U
541-73-1-----	1,3-Dichlorobenzene	18000	U
106-46-7-----	1,4-Dichlorobenzene	18000	U
100-51-6-----	Benzyl Alcohol	18000	U
95-50-1-----	1,2-Dichlorobenzene	18000	U
95-48-7-----	2-Methylphenol	150000	D
39638-32-9-----	bis(2-Chloroisopropyl) Ether	18000	U
106-44-5-----	4-Methylphenol	240000	D
621-64-7-----	N-Nitroso-Di-n-Propylamine	18000	U
67-72-1-----	Hexachloroethane	18000	U
98-95-3-----	Nitrobenzene	18000	U
78-59-1-----	Isophorone	18000	U
88-75-5-----	2-Nitrophenol	18000	U
105-67-9-----	2,4-Dimethylphenol	220000	D
65-85-0-----	Benzoic Acid	88000	U
111-91-1-----	bis(2-Chloroethoxy) Methane	18000	U
120-83-2-----	2,4-Dichlorophenol	18000	U
120-82-1-----	1,2,4-Trichlorobenzene	18000	U
91-20-3-----	Naphthalene	57000	D
106-47-8-----	4-Chloroaniline	18000	U
87-68-3-----	Hexachlorobutadiene	18000	U
59-50-7-----	4-Chloro-3-Methylphenol	18000	U
91-57-6-----	2-Methylnaphthalene	37000	D
77-47-4-----	Hexachlorocyclopentadiene	18000	U
88-06-2-----	2,4,6-Trichlorophenol	18000	U
95-95-4-----	2,4,5-Trichlorophenol	88000	U
91-58-7-----	2-Chloronaphthalene	18000	U
88-74-4-----	2-Nitroaniline	88000	U
131-11-3-----	Dimethyl Phthalate	18000	U
208-96-8-----	Acenaphthylene	4700	DJ
606-20-2-----	2,6-Dinitrotoluene	18000	U

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201ADLMS

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309683
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GD009683C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 50

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline_____	88000	U
83-32-9-----	Acenaphthene_____	18000	U
51-28-5-----	2,4-Dinitrophenol_____	88000	U
100-02-7-----	4-Nitrophenol_____	88000	U
132-64-9-----	Dibenzofuran_____	10000	DJ
121-14-2-----	2,4-Dinitrotoluene_____	18000	U
84-66-2-----	Diethylphthalate_____	18000	U
7005-72-3-----	4-Chlorophenyl-phenylether_____	18000	U
86-73-7-----	Fluorene_____	3300	DJ
100-01-6-----	4-Nitroaniline_____	88000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol_____	88000	U
86-30-6-----	N-Nitrosodiphenylamine (1)_____	18000	U
101-55-3-----	4-Bromophenyl-phenylether_____	18000	U
118-74-1-----	Hexachlorobenzene_____	18000	U
87-86-5-----	Pentachlorophenol_____	88000	U
85-01-8-----	Phenanthrene_____	10000	DJ
120-12-7-----	Anthracene_____	3800	DJ
84-74-2-----	Di-n-Butylphthalate_____	2100	DJ
206-44-0-----	Fluoranthene_____	3000	DJ
129-00-0-----	Pyrene_____	18000	U
85-68-7-----	Butylbenzylphthalate_____	18000	U
91-94-1-----	3,3'-Dichlorobenzidine_____	36000	U
56-55-3-----	Benzo(a)Anthracene_____	18000	U
218-01-9-----	Chrysene_____	18000	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate_____	18000	U
117-84-0-----	Di-n-Octyl Phthalate_____	18000	U
205-99-2-----	Benzo(b) Fluoranthene_____	18000	U
207-08-9-----	Benzo(k) Fluoranthene_____	18000	U
50-32-8-----	Benzo(a) Pyrene_____	18000	U
193-39-5-----	Indeno(1,2,3-cd) Pyrene_____	18000	U
53-70-3-----	Dibenzo(a,h) Anthracene_____	18000	U
191-24-2-----	Benzo(g,h,i) Perylene_____	18000	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201ADLMSD

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309684
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GD009684C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 50

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2-----	Phenol	18000	U
111-44-4-----	bis(2-Chloroethyl)Ether	18000	U
95-57-8-----	2-Chlorophenol	18000	U
541-73-1-----	1,3-Dichlorobenzene	18000	U
106-46-7-----	1,4-Dichlorobenzene	18000	U
100-51-6-----	Benzyl Alcohol	18000	U
95-50-1-----	1,2-Dichlorobenzene	18000	U
95-48-7-----	2-Methylphenol	130000	D
39638-32-9-----	bis(2-Chloroisopropyl)Ether	18000	U
106-44-5-----	4-Methylphenol	200000	D
621-64-7-----	N-Nitroso-Di-n-Propylamine	18000	U
67-72-1-----	Hexachloroethane	18000	U
98-95-3-----	Nitrobenzene	18000	U
78-59-1-----	Isophorone	18000	U
88-75-5-----	2-Nitrophenol	18000	U
105-67-9-----	2,4-Dimethylphenol	160000	D
65-85-0-----	Benzoic Acid	88000	U
111-91-1-----	bis(2-Chloroethoxy)Methane	18000	U
120-83-2-----	2,4-Dichlorophenol	18000	U
120-82-1-----	1,2,4-Trichlorobenzene	18000	U
91-20-3-----	Naphthalene	45000	D
106-47-8-----	4-Chloroaniline	18000	U
87-68-3-----	Hexachlorobutadiene	18000	U
59-50-7-----	4-Chloro-3-Methylphenol	18000	U
91-57-6-----	2-Methylnaphthalene	30000	D
77-47-4-----	Hexachlorocyclopentadiene	18000	U
88-06-2-----	2,4,6-Trichlorophenol	18000	U
95-95-4-----	2,4,5-Trichlorophenol	88000	U
91-58-7-----	2-Chloronaphthalene	18000	U
88-74-4-----	2-Nitroaniline	88000	U
131-11-3-----	Dimethyl Phthalate	18000	U
208-96-8-----	Acenaphthylene	3500	DJ
606-20-2-----	2,6-Dinitrotoluene	18000	U

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201ADLMSD

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309684
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GD009684C20
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: 6.0 Dilution Factor: 50

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2	3-Nitroaniline	88000	U
83-32-9	Acenaphthene	18000	U
51-28-5	2,4-Dinitrophenol	88000	U
100-02-7	4-Nitrophenol	88000	U
132-64-9	Dibenzofuran	7500	DJ
121-14-2	2,4-Dinitrotoluene	18000	U
84-66-2	Diethylphthalate	18000	U
7005-72-3	4-Chlorophenyl-phenylether	18000	U
86-73-7	Fluorene	2100	DJ
100-01-6	4-Nitroaniline	88000	U
534-52-1	4,6-Dinitro-2-Methylphenol	88000	U
86-30-6	N-Nitrosodiphenylamine (1)	18000	U
101-55-3	4-Bromophenyl-phenylether	18000	U
118-74-1	Hexachlorobenzene	18000	U
87-86-5	Pentachlorophenol	88000	U
85-01-8	Phenanthrene	16000	DJ
120-12-7	Anthracene	2200	DJ
84-74-2	Di-n-Butylphthalate	18000	U
206-44-0	Fluoranthene	2500	DJ
129-00-0	Pyrene	18000	U
85-68-7	Butylbenzylphthalate	18000	U
91-94-1	3,3'-Dichlorobenzidine	36000	U
56-55-3	Benzo(a)Anthracene	18000	U
218-01-9	Chrysene	18000	U
117-81-7	bis(2-Ethylhexyl) Phthalate	18000	U
117-84-0	Di-n-Octyl Phthalate	18000	U
205-99-2	Benzo(b) Fluoranthene	18000	U
207-08-9	Benzo(k) Fluoranthene	18000	U
50-32-8	Benzo(a) Pyrene	18000	U
193-39-5	Indeno(1,2,3-cd) Pyrene	18000	U
53-70-3	Dibenzo(a,h) Anthracene	18000	U
191-24-2	Benzo(g,h,i) Perylene	18000	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TARMS

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 310767
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GRD10767A08
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: 8.5 Dilution Factor: 5.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	280000	U
111-44-4	bis(2-Chloroethyl) Ether	280000	U
95-57-8	2-Chlorophenol	280000	U
541-73-1	1,3-Dichlorobenzene	280000	U
106-46-7	1,4-Dichlorobenzene	280000	U
100-51-6	Benzyl Alcohol	280000	U
95-50-1	1,2-Dichlorobenzene	280000	U
95-48-7	2-Methylphenol	1200000	
39638-32-9	bis(2-Chloroisopropyl) Ether	280000	U
106-44-5	4-Methylphenol	1500000	
621-64-7	N-Nitroso-Di-n-Propylamine	280000	U
67-72-1	Hexachloroethane	280000	U
98-95-3	Nitrobenzene	280000	U
78-59-1	Isophorone	280000	U
88-75-5	2-Nitrophenol	280000	U
105-67-9	2,4-Dimethylphenol	2000000	
65-85-0	Benzoic Acid	1300000	U
111-91-1	bis(2-Chloroethoxy) Methane	280000	U
120-83-2	2,4-Dichlorophenol	280000	U
120-82-1	1,2,4-Trichlorobenzene	280000	U
91-20-3	Naphthalene	320000	
106-47-8	4-Chloroaniline	280000	U
87-68-3	Hexachlorobutadiene	280000	U
59-50-7	4-Chloro-3-Methylphenol	280000	U
91-57-6	2-Methylnaphthalene	540000	
77-47-4	Hexachlorocyclopentadiene	280000	U
88-06-2	2,4,6-Trichlorophenol	280000	U
95-95-4	2,4,5-Trichlorophenol	1300000	U
91-58-7	2-Chloronaphthalene	280000	U
88-74-4	2-Nitroaniline	1300000	U
131-11-3	Dimethyl Phthalate	280000	U
208-96-8	Acenaphthylene	280000	U
606-20-2	2,6-Dinitrotoluene	280000	U

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TARMS

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 310767
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GRD10767A08
 Level: (low/med) MED Date Received: 12/20/89
 ‡ Moisture: not dec. 64 dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: 8.5 Dilution Factor: 5.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG Q

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline	1300000	U
83-32-9-----	Acenaphthene	280000	U
51-28-5-----	2,4-Dinitrophenol	1300000	U
100-02-7-----	4-Nitrophenol	1300000	U
132-64-9-----	Dibenzofuran	50000	J
121-14-2-----	2,4-Dinitrotoluene	280000	U
84-66-2-----	Diethylphthalate	280000	U
7005-72-3-----	4-Chlorophenyl-phenylether	280000	U
86-73-7-----	Fluorene	110000	J
100-01-6-----	4-Nitroaniline	1300000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	1300000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	280000	U
101-55-3-----	4-Bromophenyl-phenylether	280000	U
118-74-1-----	Hexachlorobenzene	280000	U
87-86-5-----	Pentachlorophenol	1300000	U
85-01-8-----	Phenanthrene	280000	U
120-12-7-----	Anthracene	280000	U
84-74-2-----	Di-n-Butylphthalate	280000	U
206-44-0-----	Fluoranthene	280000	U
129-00-0-----	Pyrene	280000	U
85-68-7-----	Butylbenzylphthalate	280000	U
91-94-1-----	3,3'-Dichlorobenzidine	550000	U
56-55-3-----	Benzo(a)Anthracene	280000	U
218-01-9-----	Chrysene	280000	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	280000	U
117-84-0-----	Di-n-Octyl Phthalate	280000	U
205-99-2-----	Benzo(b)Fluoranthene	280000	U
207-08-9-----	Benzo(k)Fluoranthene	280000	U
50-32-8-----	Benzo(a)Pyrene	280000	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	280000	U
53-70-3-----	Dibenzo(a,h)Anthracene	280000	U
191-24-2-----	Benzo(g,h,i)Perylene	280000	U

(1) - Cannot be separated from Diphenylamine

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TARMSD

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 310768
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GRD10768A08
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: 8.5 Dilution Factor: 5.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	<u>Q</u>
108-95-2	Phenol	280000	U
111-44-4	bis(2-Chloroethyl) Ether	280000	U
95-57-8	2-Chlorophenol	280000	U
541-73-1	1,3-Dichlorobenzene	280000	U
106-46-7	1,4-Dichlorobenzene	280000	U
100-51-6	Benzyl Alcohol	280000	U
95-50-1	1,2-Dichlorobenzene	280000	U
95-48-7	2-Methylphenol	1100000	
39638-32-9	bis(2-Chloroisopropyl) Ether	280000	U
106-44-5	4-Methylphenol	1400000	
621-64-7	N-Nitroso-Di-n-Propylamine	280000	U
67-72-1	Hexachloroethane	280000	U
98-95-3	Nitrobenzene	280000	U
78-59-1	Isophorone	280000	U
88-75-5	2-Nitrophenol	280000	U
105-67-9	2,4-Dimethylphenol	1800000	
65-85-0	Benzoic Acid	1300000	U
111-91-1	bis(2-Chloroethoxy) Methane	280000	U
120-83-2	2,4-Dichlorophenol	280000	U
120-82-1	1,2,4-Trichlorobenzene	280000	U
91-20-3	Naphthalene	220000	J
106-47-8	4-Chloroaniline	280000	U
87-68-3	Hexachlorobutadiene	280000	U
59-50-7	4-Chloro-3-Methylphenol	280000	U
91-57-6	2-Methylnaphthalene	410000	
77-47-4	Hexachlorocyclopentadiene	280000	U
88-06-2	2,4,6-Trichlorophenol	280000	U
95-95-4	2,4,5-Trichlorophenol	1300000	U
91-58-7	2-Chloronaphthalene	280000	U
88-74-4	2-Nitroaniline	1300000	U
131-11-3	Dimethyl Phthalate	280000	U
208-96-8	Acenaphthylene	280000	U
606-20-2	2,6-Dinitrotoluene	280000	U

FORM I SV-1

1/87 Rev.

1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TARMSD

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 310768
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GRD10768A08
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: 8.5 Dilution Factor: 5.0

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

99-09-2-----3-Nitroaniline	1300000	U
83-32-9-----Acenaphthene	280000	U
51-28-5-----2,4-Dinitrophenol	1300000	U
100-02-7-----4-Nitrophenol	1300000	U
132-64-9-----Dibenzofuran	38000	J
121-14-2-----2,4-Dinitrotoluene	280000	U
84-66-2-----Diethylphthalate	280000	U
7005-72-3-----4-Chlorophenyl-phenylether	280000	U
86-73-7-----Fluorene	75000	J
100-01-6-----4-Nitroaniline	1300000	U
534-52-1-----4,6-Dinitro-2-Methylphenol	1300000	U
86-30-6-----N-Nitrosodiphenylamine (1)	280000	U
101-55-3-----4-Bromophenyl-phenylether	280000	U
118-74-1-----Hexachlorobenzene	280000	U
87-86-5-----Pentachlorophenol	1300000	U
85-01-8-----Phenanthrene	280000	U
120-12-7-----Anthracene	280000	U
84-74-2-----Di-n-Butylphthalate	280000	U
206-44-0-----Fluoranthene	280000	U
129-00-0-----Pyrene	280000	U
85-68-7-----Butylbenzylphthalate	280000	U
91-94-1-----3,3'-Dichlorobenzidine	550000	U
56-55-3-----Benzo(a)Anthracene	280000	U
218-01-9-----Chrysene	280000	U
117-81-7-----bis(2-Ethylhexyl) Phthalate	280000	U
117-84-0-----Di-n-Octyl Phthalate	280000	U
205-99-2-----Benzo(b) Fluoranthene	280000	U
207-08-9-----Benzo(k) Fluoranthene	280000	U
50-32-8-----Benzo(a) Pyrene	280000	U
193-39-5-----Indeno(1,2,3-cd) Pyrene	280000	U
53-70-3-----Dibenzo(a,h) Anthracene	280000	U
191-24-2-----Benzo(g,h,i) Perylene	280000	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

3. By fraction (VOA, SV, PEST) - surrogate spike analysis results (Form II) by matrix (Water and/or Soil) and for soil, by concentration (Low or Medium)

2B
SOIL VOLATILE SURROGATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Level: (low/med) LOW

	EPA SAMPLE NO.	S1 (TOL) #	S2 (BFB) #	S3 (DCE) #	OTHER	TOT OUT
01	B201A	80 *	90	97		1
02	B201B	90	99	94		0
03	B202A	95	92	97		0
04	B202B	92	96	97		0
05	B202C	94	91	97		0
06	B202TAR	88	93	97		0
07	B202AMS	101	98	92		0
08	B202AMSD	114	113	102		0
09	VBLKL8	107	98	95		0
10	VBLKB4	97	87	97		0
11	VBLKF5	92	91	92		0

QC LIMITS

S1 (TOL) = Toluene-d8 (81-117)
 S2 (BFB) = Bromofluorobenzene (74-121)
 S3 (DCE) = 1,2-Dichloroethane-d4 (70-121)

Column to be used to flag recovery values
 * Values outside of contract required QC limits
 D Surrogates diluted out

2B
SOIL VOLATILE SURROGATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Level: (low/med) MED

	EPA SAMPLE NO.	S1 (TOL) #	S2 (BFB) #	S3 (DCE) #	OTHER	TOT OUT
01	B201ARE	91	93	111		0
02	B201BRE	115	105	113		0
03	B202TARDL	82	110	109		0
04	B201BREMS	90	100	90		0
05	B201BREMSD	86	101	104		0
06	VBLKH1	110	103	117		0
07	VBLKH9	101	104	117		0
08	VBLKI6	112	106	121		0

QC LIMITS

S1 (TOL) = Toluene-d8 (81-117)
 S2 (BFB) = Bromofluorobenzene (74-121)
 S3 (DCE) = 1,2-Dichloroethane-d4 (70-121)

Column to be used to flag recovery values
 * Values outside of contract required QC limits
 D Surrogates diluted out

2D
SOIL SEMIVOLATILE SURROGATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Level: (low/med) LOW

	EPA SAMPLE NO.	S1 (NBZ) #	S2 (FBP) #	S3 (TPH) #	S4 (PHL) #	S5 (2FP) #	S6 (TBP) #	OTHER	TOT OUT
01	B201A	556 *	97	66	70	63	35		1
02	B201ADL	D	D	D	D	D	D		0
03	B202A	63	69	84	67	71	81		0
04	B202B	66	64	77	81	79	84		0
05	B202BDL	70	69	77	59	67	77		0
06	B202C	86	86	108	86	88	98		0
07	B201ADLMS	D	D	D	D	D	D		0
08	B201ADLMSD	D	D	D	D	D	D		0
09	SBLK44	67	65	85	71	73	48		0
10	SBLK44	87	100	106	78	62	38		0

QC LIMITS

S1 (NBZ) = Nitrobenzene-d5 (23-120)
 S2 (FBP) = 2-Fluorobiphenyl (30-115)
 S3 (TPH) = Terphenyl (18-137)
 S4 (PHL) = Phenol-d5 (24-113)
 S5 (2FP) = 2-Fluorophenol (25-121)
 S6 (TBP) = 2,4,6-Tribromophenol (19-122)

Column to be used to flag recovery values
 * Values outside of contract required QC limits
 D Surrogates diluted out

2D
SOIL SEMIVOLATILE SURROGATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Level: (low/med) MED

	EPA SAMPLE NO.	S1 (NBZ) #	S2 (FBP) #	S3 (TPH) #	S4 (PHL) #	S5 (2FP) #	S6 (TBP) #	OTHER	TOT OUT
01	B201B	90	102	110	83	73	84		0
02	B202TAR	86	102	100	59	53	31		0
03	B202TARMS	100	112	105	72	71	64		0
04	B202TARMSD	77	100	94	78	73	105		0
05	SBLK82	55	81	78	71	69	73		0
06	SBLK79	57	73	72	44	47	63		0

QC LIMITS

S1 (NBZ) = Nitrobenzene-d5 (23-120)
 S2 (FBP) = 2-Fluorobiphenyl (30-115)
 S3 (TPH) = Terphenyl (18-137)
 S4 (PHL) = Phenol-d5 (24-113)
 S5 (2FP) = 2-Fluorophenol (25-121)
 S6 (TBP) = 2,4,6-Tribromophenol (19-122)

Column to be used to flag recovery values
 * Values outside of contract required QC limits
 D Surrogates diluted out

4. By fraction (VOA, SV, PEST) - matrix spike duplicate results
(Form III)

SOIL VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM LABSContract: (2-88)-REVSLab Code: COMPUCase No.: 18756

SAS No.: _____

SDG No.: 07Matrix Spike - EPA Sample No.: B202ALevel: (low/med) LOW

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
1,1-Dichloroethene	56.8	0	58.1	102	59-172
Trichloroethene	56.8	0	47.3	83	62-137
Benzene	56.8	0	55.2	97	66-142
Toluene	56.8	1.45	53.9	92	59-139
Chlorobenzene	56.8	0	52.0	92	60-133

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
1,1-Dichloroethene	56.8	53.5	94	8	22	59-172
Trichloroethene	56.8	47.5	84	-1	24	62-137
Benzene	56.8	55.1	97	0	21	66-142
Toluene	56.8	54.0	93	-1	21	59-139
Chlorobenzene	56.8	50.9	90	2	21	60-133

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 0 out of 5 outside limitsSpike Recovery: 0 out of 10 outside limitsCOMMENTS: CLP ,1875,67, , ,SOIL,309688,VOLATILE, ,
TUNE: 0019 122289 0013

SOIL VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix Spike - EPA Sample No.: B201BRE Level: (low/med) MED

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
1,1-Dichloroethene	7100	0	4730	67	59-172
Trichloroethene	7100	0	6950	98	62-137
Benzene	7100	2160	9740	107	66-142
Toluene	7100	15900	18900	42 *	59-139
Chlorobenzene	7100	0	6970	98	60-133

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS	
					RPD	REC.
1,1-Dichloroethene	7100	4330	61	9	22	59-172
Trichloroethene	7100	6970	98	0	24	62-137
Benzene	7100	10500	118	-10	21	66-142
Toluene	7100	18000	30 *	33 *	21	59-139
Chlorobenzene	7100	7230	102	-4	21	60-133

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 1 out of 5 outside limits

Spike Recovery: 2 out of 10 outside limits

COMMENTS: CLP ,1875,67, , ,SOIL,309686,VOLATILE, ,
TUNE: 0013 122989 1535

3D
SOIL SEMIVOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix Spike - EPA Sample No.: B201ADL Level: (low/med) LOW

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
Phenol	7340	22900	0	-312 *	26- 90
2-Chlorophenol	7340	0	0	0 *	25-102
1,4-Dichlorobenzene	3670	0	0	0 *	28 104
N-Nitroso-di-n-prop. (1)	3670	0	0	0 *	41 126
1,2,4-Trichlorobenzene	3670	0	0	0 *	38 107
4-Chloro-3-methylphenol	7340	0	0	0 *	26 103
Acenaphthene	3670	2590	0	-71 *	31-137
4-Nitrophenol	7340	0	0	0 *	11-114
2,4-Dinitrotoluene	3670	0	0	0 *	28- 89
Pentachlorophenol	7340	0	0	0 *	17-109
Pyrene	3670	2920	0	-80 *	35-142

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD REC.
Phenol	7340	0	-312 *	0	35 26- 90
2-Chlorophenol	7340	0	0 *	0	50 25-102
1,4-Dichlorobenzene	3670	0	0 *	0	27 28 104
N-Nitroso-di-n-prop. (1)	3670	0	0 *	0	38 41 126
1,2,4-Trichlorobenzene	3670	0	0 *	0	23 38 107
4-Chloro-3-methylphenol	7340	0	0 *	0	33 26 103
Acenaphthene	3670	0	-71 *	0	19 31-137
4-Nitrophenol	7340	0	0 *	0	50 11-114
2,4-Dinitrotoluene	3670	0	0 *	0	47 28- 89
Pentachlorophenol	7340	0	0 *	0	47 17-109
Pyrene	3670	0	-80 *	0	36 35-142

(1) N-Nitroso-di-n-propylamine

Column to be used to flag recovery and RPD values with an asterisk
 * Values outside of QC limits

RPD: 0 out of 11 outside limits
 Spike Recovery: 22 out of 22 outside limits

COMMENTS: CLP ,1875,67, , ,SOIL,309679,BNA, ,
 TUNE: 0020 122289 0034

SOIL SEMIVOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM LABSContract: (2-88)-REVSLab Code: COMPUCase No.: 18756

SAS No.: _____

SDG No.: 07Matrix Spike - EPA Sample No.: B202TARLevel: (low/med) MED

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
Phenol	1110000	331000	825000	45	26- 90
2-Chlorophenol	1110000	0	458000	41	25-102
1,4-Dichlorobenzene	556000	0	216000	39	28 104
N-Nitroso-di-n-prop. (1)	556000	0	225000	40 *	41 126
1,2,4-Trichlorobenzene	556000	0	201000	36 *	38 107
4-Chloro-3-methylphenol	1110000	0	675000	61	26 103
Acenaphthene	556000	0	431000	78	31-137
4-Nitrophenol	1110000	0	389000	35	11-114
2,4-Dinitrotoluene	556000	0	277000	50	28- 89
Pentachlorophenol	1110000	0	197000	18	17-109
Pyrene	556000	0	264000	47	35-142

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
Phenol	1110000	836000	46	-2	35	26- 90
2-Chlorophenol	1110000	464000	42	-2	50	25-102
1,4-Dichlorobenzene	556000	124000	22 *	56 *	27	28 104
N-Nitroso-di-n-prop. (1)	556000	181000	33 *	19	38	41 126
1,2,4-Trichlorobenzene	556000	157000	28 *	25 *	23	38 107
4-Chloro-3-methylphenol	1110000	658000	59	3	33	26 103
Acenaphthene	556000	353000	63	21 *	19	31-137
4-Nitrophenol	1110000	400000	36	-3	50	11-114
2,4-Dinitrotoluene	556000	276000	50	0	47	28- 89
Pentachlorophenol	1110000	191000	17	6	47	17-109
Pyrene	556000	253000	46	2	36	35-142

(1) N-Nitroso-di-n-propylamine

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 3 out of 11 outside limitsSpike Recovery: 5 out of 22 outside limitsCOMMENTS: CLP ,1875,67, , ,SOIL,309687,BNA, ,
TUNE: 0008 011090 0610

5. By fraction (VOA, SV, PEST) - blank data (Form IV) and tabulated results (Form I) including tentatively identified compounds (Form I, TIC) (VOA and SV only)

4A
VOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: GH009785C19 Lab Sample ID: VBLKB4
 Date Analyzed: 12/22/89 Time Analyzed: 0156
 Matrix: (soil/water) SOIL Level: (low/med) LOW
 Instrument ID: 19

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	B201A	309679	GH009679C19	0418
02	B202A	309688	GH009688C19	0656
03	B202B	309689	GH009689C19	0740
04	B202C	309690	GH009690A19	0815

COMMENTS: CLP , , , , , 309785, VOLATILE, BLANK,
TUNE: 0019 122289 0013

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLKB4

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKB4
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009785C19
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ Date Analyzed: 12/22/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	Chloromethane	10	U
74-83-9	Bromomethane	10	U
75-01-4	Vinyl Chloride	10	U
75-00-3	Chloroethane	10	U
75-09-2	Methylene Chloride	9	
67-64-1	Acetone	5	J
75-15-0	Carbon Disulfide	5	U
75-35-4	1,1-Dichloroethene	5	U
75-34-3	1,1-Dichloroethane	5	U
540-59-0	1,2-Dichloroethene (total)	5	U
67-66-3	Chloroform	5	U
107-06-2	1,2-Dichloroethane	5	U
78-93-3	2-Butanone	10	U
71-55-6	1,1,1-Trichloroethane	5	U
56-23-5	Carbon Tetrachloride	5	U
108-05-4	Vinyl Acetate	10	U
75-27-4	Bromodichloromethane	5	U
78-87-5	1,2-Dichloropropane	5	U
10061-01-5	cis-1,3-Dichloropropene	5	U
79-01-6	Trichloroethene	5	U
124-48-1	Dibromochloromethane	5	U
79-00-5	1,1,2-Trichloroethane	5	U
71-43-2	Benzene	5	U
10061-02-6	Trans-1,3-Dichloropropene	5	U
75-25-2	Bromoform	5	U
108-10-1	4-Methyl-2-Pentanone	10	U
591-78-6	2-Hexanone	10	U
127-18-4	Tetrachloroethene	5	U
79-34-5	1,1,2,2-Tetrachloroethane	5	U
108-88-3	Toluene	5	U
108-90-7	Chlorobenzene	5	U
100-41-4	Ethylbenzene	5	U
100-42-5	Styrene	5	U
1330-20-7	Total Xylenes	5	U

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1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLKB4

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKB4
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009785C19
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

4A
VOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: GH010660C19 Lab Sample ID: VBLKF5
 Date Analyzed: 12/27/89 Time Analyzed: 0229
 Matrix: (soil/water) SOIL Level: (low/med) LOW
 Instrument ID: 19

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	B201B	309686	GR009686C19	0539
02	B202TAR	309687	GR009687C19	0710

COMMENTS: CLP , , , , , 310660, VOLATILE, BLANK,
 TUNE: 0019 122789 0032

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EPA SAMPLE NO.

VBLKF5

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKF5
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH010660C19
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ Date Analyzed: 12/27/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	10	U
74-83-9	-----Bromomethane	10	U
75-01-4	-----Vinyl Chloride	10	U
75-00-3	-----Chloroethane	10	U
75-09-2	-----Methylene Chloride	3	J
67-64-1	-----Acetone	3	J
75-15-0	-----Carbon Disulfide	5	U
75-35-4	-----1,1-Dichloroethene	5	U
75-34-3	-----1,1-Dichloroethane	5	U
540-59-0	-----1,2-Dichloroethene (total)	5	U
67-66-3	-----Chloroform	5	U
107-06-2	-----1,2-Dichloroethane	5	U
78-93-3	-----2-Butanone	10	U
71-55-6	-----1,1,1-Trichloroethane	5	U
56-23-5	-----Carbon Tetrachloride	5	U
108-05-4	-----Vinyl Acetate	10	U
75-27-4	-----Bromodichloromethane	5	U
78-87-5	-----1,2-Dichloropropane	5	U
10061-01-5	-----cis-1,3-Dichloropropene	5	U
79-01-6	-----Trichloroethene	5	U
124-48-1	-----Dibromochloromethane	5	U
79-00-5	-----1,1,2-Trichloroethane	5	U
71-43-2	-----Benzene	5	U
10061-02-6	-----Trans-1,3-Dichloropropene	5	U
75-25-2	-----Bromoform	5	U
108-10-1	-----4-Methyl-2-Pentanone	10	U
591-78-6	-----2-Hexanone	3	J
127-18-4	-----Tetrachloroethene	5	U
79-34-5	-----1,1,2,2-Tetrachloroethane	5	U
108-88-3	-----Toluene	5	U
108-90-7	-----Chlorobenzene	5	U
100-41-4	-----Ethylbenzene	5	U
100-42-5	-----Styrene	5	U
1330-20-7	-----Total Xylenes	5	U

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 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLKF5

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKF5
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH010660C19
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ Date Analyzed: 12/27/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

4A
VOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: CN011246B13 Lab Sample ID: VBLKH1
 Date Analyzed: 12/28/89 Time Analyzed: 1742
 Matrix: (soil/water) SOIL Level: (low/med) MED
 Instrument ID: 13

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01 B202TARDL	309687	C3R09687C13	0152

COMMENTS: CLP , , , , , 311246, VOLATILE, BLANK,
 TUNE: 0013 122889 1528

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EPA SAMPLE NO.

VBLKH1

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKH1
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: CN011246B13
 Level: (low/med) MED Date Received: _____
 ‡ Moisture: not dec. _____ Date Analyzed: 12/28/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	1300	U
74-83-9	-----Bromomethane	1300	U
75-01-4	-----Vinyl Chloride	1300	U
75-00-3	-----Chloroethane	1300	U
75-09-2	-----Methylene Chloride	290	J
67-64-1	-----Acetone	1300	U
75-15-0	-----Carbon Disulfide	630	U
75-35-4	-----1,1-Dichloroethene	630	U
75-34-3	-----1,1-Dichloroethane	630	U
540-59-0	-----1,2-Dichloroethene (total)	630	U
67-66-3	-----Chloroform	630	U
107-06-2	-----1,2-Dichloroethane	630	U
78-93-3	-----2-Butanone	1300	U
71-55-6	-----1,1,1-Trichloroethane	630	U
56-23-5	-----Carbon Tetrachloride	630	U
108-05-4	-----Vinyl Acetate	1300	U
75-27-4	-----Bromodichloromethane	630	U
78-87-5	-----1,2-Dichloropropane	630	U
10061-01-5	-----cis-1,3-Dichloropropene	630	U
79-01-6	-----Trichloroethene	630	U
124-48-1	-----Dibromochloromethane	630	U
79-00-5	-----1,1,2-Trichloroethane	630	U
71-43-2	-----Benzene	630	U
10061-02-6	-----Trans-1,3-Dichloropropene	630	U
75-25-2	-----Bromoform	630	U
108-10-1	-----4-Methyl-2-Pentanone	1300	U
591-78-6	-----2-Hexanone	1300	U
127-18-4	-----Tetrachloroethene	630	U
79-34-5	-----1,1,2,2-Tetrachloroethane	630	U
108-88-3	-----Toluene	630	U
108-90-7	-----Chlorobenzene	630	U
100-41-4	-----Ethylbenzene	630	U
100-42-5	-----Styrene	630	U
1330-20-7	-----Total Xylenes	630	U

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 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLKH1

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKH1
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: CN011246B13
 Level: (low/med) MED Date Received: _____
 ‡ Moisture: not dec. _____ Date Analyzed: 12/28/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

4A
VOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: CN011397B13 Lab Sample ID: VBLKH9
 Date Analyzed: 12/28/89 Time Analyzed: 1900
 Matrix: (soil/water) SOIL Level: (low/med) MED
 Instrument ID: 13

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	B201ARE	309679	C3R09679B13	2034

COMMENTS: CLP , , , , , 311397, VOLATILE, BLANK,
 TUNE: 0013 122889 1528

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLKH9

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKH9
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: CN011397B13
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ Date Analyzed: 12/28/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	1300	U
74-83-9	-----Bromomethane	1300	U
75-01-4	-----Vinyl Chloride	1300	U
75-00-3	-----Chloroethane	1300	U
75-09-2	-----Methylene Chloride	290	J
67-64-1	-----Acetone	1300	U
75-15-0	-----Carbon Disulfide	630	U
75-35-4	-----1,1-Dichloroethene	630	U
75-34-3	-----1,1-Dichloroethane	630	U
540-59-0	-----1,2-Dichloroethene (total)	630	U
67-66-3	-----Chloroform	630	U
107-06-2	-----1,2-Dichloroethane	630	U
78-93-3	-----2-Butanone	1300	U
71-55-6	-----1,1,1-Trichloroethane	630	U
56-23-5	-----Carbon Tetrachloride	630	U
108-05-4	-----Vinyl Acetate	1300	U
75-27-4	-----Bromodichloromethane	630	U
78-87-5	-----1,2-Dichloropropane	630	U
10061-01-5	-----cis-1,3-Dichloropropene	630	U
79-01-6	-----Trichloroethene	630	U
124-48-1	-----Dibromochloromethane	630	U
79-00-5	-----1,1,2-Trichloroethane	630	U
71-43-2	-----Benzene	630	U
10061-02-6	-----Trans-1,3-Dichloropropene	630	U
75-25-2	-----Bromoform	630	U
108-10-1	-----4-Methyl-2-Pentanone	1300	U
591-78-6	-----2-Hexanone	1300	U
127-18-4	-----Tetrachloroethene	630	U
79-34-5	-----1,1,2,2-Tetrachloroethane	630	U
108-88-3	-----Toluene	630	U
108-90-7	-----Chlorobenzene	630	U
100-41-4	-----Ethylbenzene	630	U
100-42-5	-----Styrene	630	U
1330-20-7	-----Total Xylenes	630	U

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1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLKH9

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKH9
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: CN011397B13
 Level: (low/med) MED Date Received: _____
 ‡ Moisture: not dec. _____ Date Analyzed: 12/28/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

4A
VOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: CR011532B13 Lab Sample ID: VELKI6
 Date Analyzed: 12/29/89 Time Analyzed: 1829
 Matrix: (soil/water): SOIL Level: (low/med) MED
 Instrument ID: 13

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	B201BRE	309686	C5R09686B13	2010
02	B201BREMS	311271	CN011271B13	2154
03	B201BREMSD	311272	CN011272B13	2234

COMMENTS: CLP , , , , , 311532, VOLATILE, BLANK,
 TUNE: 0013 122989 1535

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLKI6

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKI6
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: CR011532B13
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ Date Analyzed: 12/29/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	1300	U
74-83-9	-----Bromomethane	1300	U
75-01-4	-----Vinyl Chloride	1300	U
75-00-3	-----Chloroethane	1300	U
75-09-2	-----Methylene Chloride	140	J
67-64-1	-----Acetone	1300	U
75-15-0	-----Carbon Disulfide	630	U
75-35-4	-----1,1-Dichloroethene	630	U
75-34-3	-----1,1-Dichloroethane	630	U
540-59-0	-----1,2-Dichloroethene (total)	630	U
67-66-3	-----Chloroform	630	U
107-06-2	-----1,2-Dichloroethane	630	U
78-93-3	-----2-Butanone	1300	U
71-55-6	-----1,1,1-Trichloroethane	630	U
56-23-5	-----Carbon Tetrachloride	630	U
108-05-4	-----Vinyl Acetate	1300	U
75-27-4	-----Bromodichloromethane	630	U
78-87-5	-----1,2-Dichloropropane	630	U
10061-01-5	-----cis-1,3-Dichloropropene	630	U
79-01-6	-----Trichloroethene	630	U
124-48-1	-----Dibromochloromethane	630	U
79-00-5	-----1,1,2-Trichloroethane	630	U
71-43-2	-----Benzene	630	U
10061-02-6	-----Trans-1,3-Dichloropropene	630	U
75-25-2	-----Bromoform	630	U
108-10-1	-----4-Methyl-2-Pentanone	1300	U
591-78-6	-----2-Hexanone	1300	U
127-18-4	-----Tetrachloroethene	630	U
79-34-5	-----1,1,2,2-Tetrachloroethane	630	U
108-88-3	-----Toluene	630	U
108-90-7	-----Chlorobenzene	630	U
100-41-4	-----Ethylbenzene	630	U
100-42-5	-----Styrene	630	U
1330-20-7	-----Total Xylenes	630	U

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 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLKI6

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKI6
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: CR011532B13
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ Date Analyzed: 12/29/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=====	=====	=====	=====	=====

4A
VOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: GH011975A19 Lab Sample ID: VBLKL8
 Date Analyzed: 01/02/90 Time Analyzed: 1119
 Matrix: (soil/water) SOIL Level: (low/med) LOW
 Instrument ID: 19

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	B202AMS	309680	G3R09680A19	1352
02	B202AMSD	309681	G3R09681A19	1426

COMMENTS: CLP, , , , , 311975, VOLATILE, BLANK,
 TUNE: 0019 010290 0924

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VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

VBLKL8

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: VBLKL8

Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH011975A19

Level: (low/med) LOW Date Received: _____

‡ Moisture: not dec. _____ Date Analyzed: 01/02/90

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	10	U
74-83-9	-----Bromomethane	10	U
75-01-4	-----Vinyl Chloride	10	U
75-00-3	-----Chloroethane	10	U
75-09-2	-----Methylene Chloride	5	
67-64-1	-----Acetone	10	U
75-15-0	-----Carbon Disulfide	5	U
75-35-4	-----1,1-Dichloroethene	5	U
75-34-3	-----1,1-Dichloroethane	5	U
540-59-0	-----1,2-Dichloroethene (total)	5	U
67-66-3	-----Chloroform	5	U
107-06-2	-----1,2-Dichloroethane	5	U
78-93-3	-----2-Butanone	10	U
71-55-6	-----1,1,1-Trichloroethane	5	U
56-23-5	-----Carbon Tetrachloride	5	U
108-05-4	-----Vinyl Acetate	10	U
75-27-4	-----Bromodichloromethane	5	U
78-87-5	-----1,2-Dichloropropane	5	U
10061-01-5	-----cis-1,3-Dichloropropene	5	U
79-01-6	-----Trichloroethene	5	U
124-48-1	-----Dibromochloromethane	5	U
79-00-5	-----1,1,2-Trichloroethane	5	U
71-43-2	-----Benzene	5	U
10061-02-6	-----Trans-1,3-Dichloropropene	5	U
75-25-2	-----Bromoform	5	U
108-10-1	-----4-Methyl-2-Pentanone	10	U
591-78-6	-----2-Hexanone	10	U
127-18-4	-----Tetrachloroethene	5	U
79-34-5	-----1,1,2,2-Tetrachloroethane	5	U
108-88-3	-----Toluene	5	U
108-90-7	-----Chlorobenzene	5	U
100-41-4	-----Ethylbenzene	5	U
100-42-5	-----Styrene	5	U
1330-20-7	-----Total Xylenes	5	U

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1E
 VOLATILE ORGANICS ANALYSIS DATA SHEET
 TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

VBLKL8

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: VBLKL8
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH011975A19
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ Date Analyzed: 01/02/90
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS:
 (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

4B
SEMIVOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: G2J10176C20 Lab Sample ID: SBLK44
 Date Extracted: 12/21/89 Extraction: (SepF/Cont/Sonc) SONC
 Date Analyzed: 12/22/89 Time Analyzed: 0034
 Matrix: (soil/water) SOIL Level: (low/med) LOW
 Instrument ID: 20

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
01	B201ADL	309679	G2D09679C20	12/22/89
02	B202A	309688	GH009688C20	12/22/89
03	B202B	309689	GH009689C20	12/22/89
04	B202BDL	309689	GD009689A20	12/22/89
05	B202C	309690	GH009690A20	12/22/89
06	B201ADLMS	309683	GD009683C20	12/22/89
07	B201ADLMSD	309684	GD009684C20	12/22/89

COMMENTS: CLP , , , , , 310176, BNA, BLANK,
TUNE: 0020 122289 0034

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

SBLK44

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK44
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G2J10176C20
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	330	U
111-44-4	bis(2-Chloroethyl) Ether	330	U
95-57-8	2-Chlorophenol	330	U
541-73-1	1,3-Dichlorobenzene	330	U
106-46-7	1,4-Dichlorobenzene	330	U
100-51-6	Benzyl Alcohol	330	U
95-50-1	1,2-Dichlorobenzene	330	U
95-48-7	2-Methylphenol	330	U
39638-32-9	bis(2-Chloroisopropyl) Ether	330	U
106-44-5	4-Methylphenol	330	U
621-64-7	N-Nitroso-Di-n-Propylamine	330	U
67-72-1	Hexachloroethane	330	U
98-95-3	Nitrobenzene	330	U
78-59-1	Isophorone	330	U
88-75-5	2-Nitrophenol	330	U
105-67-9	2,4-Dimethylphenol	330	U
65-85-0	Benzoic Acid	1600	U
111-91-1	bis(2-Chloroethoxy) Methane	330	U
120-83-2	2,4-Dichlorophenol	330	U
120-82-1	1,2,4-Trichlorobenzene	330	U
91-20-3	Naphthalene	330	U
106-47-8	4-Chloroaniline	330	U
87-68-3	Hexachlorobutadiene	330	U
59-50-7	4-Chloro-3-Methylphenol	330	U
91-57-6	2-Methylnaphthalene	330	U
77-47-4	Hexachlorocyclopentadiene	330	U
88-06-2	2,4,6-Trichlorophenol	330	U
95-95-4	2,4,5-Trichlorophenol	1600	U
91-58-7	2-Chloronaphthalene	330	U
88-74-4	2-Nitroaniline	1600	U
131-11-3	Dimethyl Phthalate	330	U
208-96-8	Acenaphthylene	330	U
606-20-2	2,6-Dinitrotoluene	330	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

SBLK44

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK44
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G2J10176C20
 Level: (low/med) LOW Date Received: _____
 ‡ Moisture: not dec. _____ dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.00

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/Kg)	UG/KG
99-09-2	3-Nitroaniline	1600	U
83-32-9	Acenaphthene	330	U
51-28-5	2,4-Dinitrophenol	1600	U
100-02-7	4-Nitrophenol	1600	U
132-64-9	Dibenzofuran	330	U
121-14-2	2,4-Dinitrotoluene	330	U
84-66-2	Diethylphthalate	330	U
7005-72-3	4-Chlorophenyl-phenylether	330	U
86-73-7	Fluorene	330	U
100-01-6	4-Nitroaniline	1600	U
534-52-1	4,6-Dinitro-2-Methylphenol	1600	U
86-30-6	N-Nitrosodiphenylamine (1)	330	U
101-55-3	4-Bromophenyl-phenylether	330	U
118-74-1	Hexachlorobenzene	330	U
87-86-5	Pentachlorophenol	1600	U
85-01-8	Phenanthrene	330	U
120-12-7	Anthracene	330	U
84-74-2	Di-n-Butylphthalate	330	U
206-44-0	Fluoranthene	330	U
129-00-0	Pyrene	330	U
85-68-7	Butylbenzylphthalate	330	U
91-94-1	3,3'-Dichlorobenzidine	660	U
56-55-3	Benzo(a)Anthracene	330	U
218-01-9	Chrysene	330	U
117-81-7	bis(2-Ethylhexyl)Phthalate	330	U
117-84-0	Di-n-Octyl Phthalate	330	U
205-99-2	Benzo(b)Fluoranthene	330	U
207-08-9	Benzo(k)Fluoranthene	330	U
50-32-8	Benzo(a)Pyrene	330	U
193-39-5	Indeno(1,2,3-cd)Pyrene	330	U
53-70-3	Dibenzo(a,h)Anthracene	330	U
191-24-2	Benzo(g,h,i)Perylene	330	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

SBLK44

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK44
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: G2J10176C20
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/22/89
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.00

Number TICs found: 1 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	TRIBROMOPHENOL ISOMER	11.92	230	J

4B
SEMIVOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: GH010176B08 Lab Sample ID: SBLK44
 Date Extracted: 12/21/89 Extraction: (SepF/Cont/Sonc) SONC
 Date Analyzed: 12/28/89 Time Analyzed: 2041
 Matrix: (soil/water) SOIL Level: (low/med) LOW
 Instrument ID: 08

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
01	B201A	309679	G3D09679B08	12/28/89

COMMENTS: CLP , , , , , 310176, BNA, BLANK,
TUNE: 0008 122889 1109

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

SBLK44

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK44
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH010176B08
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/28/89
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.00

CONCENTRATION UNITS:
 CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

108-95-2-----Phenol	330	U
111-44-4-----bis(2-Chloroethyl) Ether	330	U
95-57-8-----2-Chlorophenol	330	U
541-73-1-----1,3-Dichlorobenzene	330	U
106-46-7-----1,4-Dichlorobenzene	330	U
100-51-6-----Benzyl Alcohol	330	U
95-50-1-----1,2-Dichlorobenzene	330	U
95-48-7-----2-Methylphenol	330	U
39638-32-9-----bis(2-Chloroisopropyl) Ether	330	U
106-44-5-----4-Methylphenol	330	U
621-64-7-----N-Nitroso-Di-n-Propylamine	330	U
67-72-1-----Hexachloroethane	330	U
98-95-3-----Nitrobenzene	330	U
78-59-1-----Isophorone	330	U
88-75-5-----2-Nitrophenol	330	U
105-67-9-----2,4-Dimethylphenol	330	U
65-85-0-----Benzoic Acid	1600	U
111-91-1-----bis(2-Chloroethoxy) Methane	330	U
120-83-2-----2,4-Dichlorophenol	330	U
120-82-1-----1,2,4-Trichlorobenzene	330	U
91-20-3-----Naphthalene	330	U
106-47-8-----4-Chloroaniline	330	U
87-68-3-----Hexachlorobutadiene	330	U
59-50-7-----4-Chloro-3-Methylphenol	330	U
91-57-6-----2-Methylnaphthalene	330	U
77-47-4-----Hexachlorocyclopentadiene	330	U
88-06-2-----2,4,6-Trichlorophenol	330	U
95-95-4-----2,4,5-Trichlorophenol	1600	U
91-58-7-----2-Chloronaphthalene	330	U
88-74-4-----2-Nitroaniline	1600	U
131-11-3-----Dimethyl Phthalate	330	U
208-96-8-----Acenaphthylene	330	U
606-20-2-----2,6-Dinitrotoluene	330	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

SBLK44

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK44
 Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH010176B08
 Level: (low/med) LOW Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/21/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/28/89
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.00

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/KG Q

99-09-2-----3-Nitroaniline	1600	U
83-32-9-----Acenaphthene	330	U
51-28-5-----2,4-Dinitrophenol	1600	U
100-02-7-----4-Nitrophenol	1600	U
132-64-9-----Dibenzofuran	330	U
121-14-2-----2,4-Dinitrotoluene	330	U
84-66-2-----Diethylphthalate	330	U
7005-72-3-----4-Chlorophenyl-phenylether	330	U
86-73-7-----Fluorene	330	U
100-01-6-----4-Nitroaniline	1600	U
534-52-1-----4,6-Dinitro-2-Methylphenol	1600	U
86-30-6-----N-Nitrosodiphenylamine (1)	330	U
101-55-3-----4-Bromophenyl-phenylether	330	U
118-74-1-----Hexachlorobenzene	330	U
87-86-5-----Pentachlorophenol	1600	U
85-01-8-----Phenanthrene	330	U
120-12-7-----Anthracene	330	U
84-74-2-----Di-n-Butylphthalate	330	U
206-44-0-----Fluoranthene	330	U
129-00-0-----Pyrene	330	U
85-68-7-----Butylbenzylphthalate	330	U
91-94-1-----3,3'-Dichlorobenzidine	660	U
56-55-3-----Benzo(a)Anthracene	330	U
218-01-9-----Chrysene	330	U
117-81-7-----bis(2-Ethylhexyl) Phthalate	63	J
117-84-0-----Di-n-Octyl Phthalate	330	U
205-99-2-----Benzo(b)Fluoranthene	330	U
207-08-9-----Benzo(k)Fluoranthene	330	U
50-32-8-----Benzo(a)Pyrene	330	U
193-39-5-----Indeno(1,2,3-cd)Pyrene	330	U
53-70-3-----Dibenzo(a,h)Anthracene	330	U
191-24-2-----Benzo(g,h,i)Perylene	330	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

SBLK44

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: SBLK44

Sample wt/vol: 30.0 (g/mL) G Lab File ID: GH010176B08

Level: (low/med) LOW Date Received: _____

% Moisture: not dec. _____ dec. _____ Date Extracted: 12/21/89

Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 12/28/89

GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.00

Number TICs found: 5 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 3074-71-3	HEPTANE, 2,3-DIMETHYL-	4.95	430	J
2. 2216-34-4	OCTANE, 4-METHYL-	5.03	400	J
3.	ALDOL	5.95	400	AJ
4.	UNKNOWN	7.97	270	J
5.	TRIBROMOPHENOL ISOMER	12.84	330	J

4B
SEMIVOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: GH011831C21 Lab Sample ID: SBLK82
 Date Extracted: 12/29/89 Extraction: (SepF/Cont/Sonc) SONC
 Date Analyzed: 01/02/90 Time Analyzed: 0723
 Matrix: (soil/water) SOIL Level: (low/med) MED
 Instrument ID: 21

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
01	B201B	309686	GD009686A21	01/02/90

COMMENTS: CLP , , , , , 311831, BNA, BLANK,
 TUNE: 0021 010290 0223

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

SBLK82

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK82
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GH011831C21
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/29/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/02/90
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	20000	U
111-44-4	bis(2-Chloroethyl) Ether	20000	U
95-57-8	2-Chlorophenol	20000	U
541-73-1	1,3-Dichlorobenzene	20000	U
106-46-7	1,4-Dichlorobenzene	20000	U
100-51-6	Benzyl Alcohol	20000	U
95-50-1	1,2-Dichlorobenzene	20000	U
95-48-7	2-Methylphenol	20000	U
39638-32-9	bis(2-Chloroisopropyl) Ether	20000	U
106-44-5	4-Methylphenol	20000	U
621-64-7	N-Nitroso-Di-n-Propylamine	20000	U
67-72-1	Hexachloroethane	20000	U
98-95-3	Nitrobenzene	20000	U
78-59-1	Isophorone	20000	U
88-75-5	2-Nitrophenol	20000	U
105-67-9	2,4-Dimethylphenol	20000	U
65-85-0	Benzoic Acid	96000	U
111-91-1	bis(2-Chloroethoxy) Methane	20000	U
120-83-2	2,4-Dichlorophenol	20000	U
120-82-1	1,2,4-Trichlorobenzene	20000	U
91-20-3	Naphthalene	20000	U
106-47-8	4-Chloroaniline	20000	U
87-68-3	Hexachlorobutadiene	20000	U
59-50-7	4-Chloro-3-Methylphenol	20000	U
91-57-6	2-Methylnaphthalene	20000	U
77-47-4	Hexachlorocyclopentadiene	20000	U
88-06-2	2,4,6-Trichlorophenol	20000	U
95-95-4	2,4,5-Trichlorophenol	96000	U
91-58-7	2-Chloronaphthalene	20000	U
88-74-4	2-Nitroaniline	96000	U
131-11-3	Dimethyl Phthalate	20000	U
208-96-8	Acenaphthylene	20000	U
606-20-2	2,6-Dinitrotoluene	20000	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

SBLK82

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK82
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GH011831C21
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/29/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/02/90
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline	96000	U
83-32-9-----	Acenaphthene	20000	U
51-28-5-----	2,4-Dinitrophenol	96000	U
100-02-7-----	4-Nitrophenol	96000	U
132-64-9-----	Dibenzofuran	20000	U
121-14-2-----	2,4-Dinitrotoluene	20000	U
84-66-2-----	Diethylphthalate	20000	U
7005-72-3-----	4-Chlorophenyl-phenylether	20000	U
86-73-7-----	Fluorene	20000	U
100-01-6-----	4-Nitroaniline	96000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	96000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	20000	U
101-55-3-----	4-Bromophenyl-phenylether	20000	U
118-74-1-----	Hexachlorobenzene	20000	U
87-86-5-----	Pentachlorophenol	96000	U
85-01-8-----	Phenanthrene	20000	U
120-12-7-----	Anthracene	20000	U
84-74-2-----	Di-n-Butylphthalate	20000	U
206-44-0-----	Fluoranthene	20000	U
129-00-0-----	Pyrene	20000	U
85-68-7-----	Butylbenzylphthalate	20000	U
91-94-1-----	3,3'-Dichlorobenzidine	40000	U
56-55-3-----	Benzo(a)Anthracene	20000	U
218-01-9-----	Chrysene	20000	U
117-81-7-----	bis(2-Ethylhexyl)Phthalate	20000	U
117-84-0-----	Di-n-Octyl Phthalate	20000	U
205-99-2-----	Benzo(b)Fluoranthene	20000	U
207-08-9-----	Benzo(k)Fluoranthene	20000	U
50-32-8-----	Benzo(a)Pyrene	20000	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	20000	U
53-70-3-----	Dibenzo(a,h)Anthracene	20000	U
191-24-2-----	Benzo(g,h,i)Perylene	20000	U

(1) - Cannot be separated from Diphenylamine

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

SBLK82

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK82
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GH011831C21
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/29/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/02/90
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

4B
SEMIVOLATILE METHOD BLANK SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID: GJ011481A08 Lab Sample ID: SBLK79
 Date Extracted: 12/28/89 Extraction: (SepF/Cont/Sonc) SONC
 Date Analyzed: 01/10/90 Time Analyzed: 1147
 Matrix: (soil/water) SOIL Level: (low/med) MED
 Instrument ID: 08

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
01	B202TAR	309687	GRD09687A08	01/10/90
02	B202TARMS	310767	GRD10767A08	01/10/90
03	B202TARMSD	310768	GRD10768A08	01/10/90

COMMENTS: CLP , , , , , 311481, BNA, BLANK,
TUNE: 0008 011090 0610

1B
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

SBLK79

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK79
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GJ011481A08
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
108-95-2	Phenol	20000	U
111-44-4	bis(2-Chloroethyl) Ether	20000	U
95-57-8	2-Chlorophenol	20000	U
541-73-1	1,3-Dichlorobenzene	20000	U
106-46-7	1,4-Dichlorobenzene	20000	U
100-51-6	Benzyl Alcohol	20000	U
95-50-1	1,2-Dichlorobenzene	20000	U
95-48-7	2-Methylphenol	20000	U
39638-32-9	bis(2-Chloroisopropyl) Ether	20000	U
106-44-5	4-Methylphenol	20000	U
621-64-7	N-Nitroso-Di-n-Propylamine	20000	U
67-72-1	Hexachloroethane	20000	U
98-95-3	Nitrobenzene	20000	U
78-59-1	Isophorone	20000	U
88-75-5	2-Nitrophenol	20000	U
105-67-9	2,4-Dimethylphenol	20000	U
65-85-0	Benzoic Acid	96000	U
111-91-1	bis(2-Chloroethoxy)Methane	20000	U
120-83-2	2,4-Dichlorophenol	20000	U
120-82-1	1,2,4-Trichlorobenzene	20000	U
91-20-3	Naphthalene	20000	U
106-47-8	4-Chloroaniline	20000	U
87-68-3	Hexachlorobutadiene	20000	U
59-50-7	4-Chloro-3-Methylphenol	20000	U
91-57-6	2-Methylnaphthalene	20000	U
77-47-4	Hexachlorocyclopentadiene	20000	U
88-06-2	2,4,6-Trichlorophenol	20000	U
95-95-4	2,4,5-Trichlorophenol	96000	U
91-58-7	2-Chloronaphthalene	20000	U
88-74-4	2-Nitroaniline	96000	U
131-11-3	Dimethyl Phthalate	20000	U
208-96-8	Acenaphthylene	20000	U
606-20-2	2,6-Dinitrotoluene	20000	U

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1C
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

SBLK79

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK79
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GJ011481A08
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
99-09-2-----	3-Nitroaniline	96000	U
83-32-9-----	Acenaphthene	20000	U
51-28-5-----	2,4-Dinitrophenol	96000	U
100-02-7-----	4-Nitrophenol	96000	U
132-64-9-----	Dibenzofuran	20000	U
121-14-2-----	2,4-Dinitrotoluene	20000	U
84-66-2-----	Diethylphthalate	20000	U
7005-72-3-----	4-Chlorophenyl-phenylether	20000	U
86-73-7-----	Fluorene	20000	U
100-01-6-----	4-Nitroaniline	96000	U
534-52-1-----	4,6-Dinitro-2-Methylphenol	96000	U
86-30-6-----	N-Nitrosodiphenylamine (1)	20000	U
101-55-3-----	4-Bromophenyl-phenylether	20000	U
118-74-1-----	Hexachlorobenzene	20000	U
87-86-5-----	Pentachlorophenol	96000	U
85-01-8-----	Phenanthrene	20000	U
120-12-7-----	Anthracene	20000	U
84-74-2-----	Di-n-Butylphthalate	20000	U
206-44-0-----	Fluoranthene	20000	U
129-00-0-----	Pyrene	20000	U
85-68-7-----	Butylbenzylphthalate	20000	U
91-94-1-----	3,3'-Dichlorobenzidine	40000	U
56-55-3-----	Benzo(a)Anthracene	20000	U
218-01-9-----	Chrysene	20000	U
117-81-7-----	bis(2-Ethylhexyl) Phthalate	20000	U
117-84-0-----	Di-n-Octyl Phthalate	20000	U
205-99-2-----	Benzo(b)Fluoranthene	20000	U
207-08-9-----	Benzo(k)Fluoranthene	20000	U
50-32-8-----	Benzo(a)Pyrene	20000	U
193-39-5-----	Indeno(1,2,3-cd)Pyrene	20000	U
53-70-3-----	Dibenzo(a,h)Anthracene	20000	U
191-24-2-----	Benzo(g,h,i)Perylene	20000	U

(1) - Cannot be separated from Diphenylamine

FORM I SV-2

1/87 Rev.

1F
SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

SBLK79

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: SBLK79
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GJ011481A08
 Level: (low/med) MED Date Received: _____
 % Moisture: not dec. _____ dec. _____ Date Extracted: 12/28/89
 Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 01/10/90
 GPC Cleanup: (Y/N) N pH: _____ Dilution Factor: 1.0

Number TICs found: 0 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q

6. By fraction (VOA, SV only) - internal standard area data (Form VIII)

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): CT891228B13 Date Analyzed: 12/28/89
 Instrument ID: 13 Time Analyzed: 1636
 Matrix: (soil/water) SOIL Level: (low/med) MED Column: (pack/cap) CAP

	IS1 (BCM) AREA #	RT	IS2 (DFB) AREA #	RT	IS3 (CBZ) AREA #	RT
12 HOUR STD	56600	5.10	191000	6.77	229000	11.42
UPPER LIMIT	113200		382000		458000	
LOWER LIMIT	28300		95500		114500	
EPA SAMPLE NO.						
01 B201ARE	49900	4.98	183000	6.60	234000	11.29
02 B202TARDL	49600	4.95	173000	6.62	250000	11.32
03 VBLKH1	52000	5.10	180000	6.75	227000	11.37
04 VBLKH9	51300	5.03	171000	6.67	225000	11.32

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): CS891229A13 Date Analyzed: 12/29/89
 Instrument ID: 13 Time Analyzed: 1555
 Matrix: (soil/water) SOIL Level: (low/med) MED Column: (pack/cap) CAP

	IS1 (BCM) AREA #	RT	IS2 (DFB) AREA #	RT	IS3 (CBZ) AREA #	RT
12 HOUR STD	59500	5.13	195000	6.82	273000	11.54
UPPER LIMIT	119000		390000		546000	
LOWER LIMIT	29750		97500		136500	
EPA SAMPLE NO.						
01 B201BRE	54500	5.10	190000	6.75	270000	11.42
02 B201BREMS	57000	5.08	163000	6.75	272000	11.42
03 B201BREMSD	51500	5.07	168000	6.75	259000	11.42
04 VBLKI6	52400	5.10	171000	6.77	269000	11.45

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): G5891222C19 Date Analyzed: 12/22/89
 Instrument ID: 19 Time Analyzed: 0013
 Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) CAP

	IS1 (BCM) AREA #	RT	IS2 (DFB) AREA #	RT	IS3 (CBZ) AREA #	RT
12 HOUR STD	44400	5.90	176000	7.70	178000	12.55
UPPER LIMIT	88800		352000		356000	
LOWER LIMIT	22200		88000		89000	
EPA SAMPLE NO.						
01 B201A	34900	5.80	134000	7.60	159000	12.50
02 B202A	49300	5.83	196000	7.63	206000	12.49
03 B202B	46800	5.80	192000	7.60	204000	12.50
04 B202C	49100	5.77	202000	7.60	211000	12.49
05 VBLKB4	49300	5.85	199000	7.65	200000	12.54

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): GT891227C19 Date Analyzed: 12/27/89
 Instrument ID: 19 Time Analyzed: 0137
 Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) CAP

	IS1 (BCM) AREA #	RT	IS2 (DFB) AREA #	RT	IS3 (CBZ) AREA #	RT
12 HOUR STD	66200	5.90	245000	7.70	250000	12.57
UPPER LIMIT	132400		490000		500000	
LOWER LIMIT	33100		122500		125000	
EPA SAMPLE NO.						
01 B201B	52500	5.80	198000	7.62	195000	12.50
02 B202TAR	68100	5.78	245000	7.60	246000	12.52
03 VBLKF5	80100	5.83	310000	7.63	311000	12.54

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): GS900102A19 Date Analyzed: 01/02/90
 Instrument ID: 19 Time Analyzed: 0942
 Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) CAP

	IS1 (BCM) AREA #	RT	IS2 (DFB) AREA #	RT	IS3 (CBZ) AREA #	RT
12 HOUR STD	52100	5.90	198000	7.68	214000	12.54
UPPER LIMIT	104200		396000		428000	
LOWER LIMIT	26050		99000		107000	
EPA SAMPLE NO.						
01 B202AMS	53700	5.88	212000	7.72	234000	12.60
02 B202AMSD	59400	5.87	223000	7.68	245000	12.60
03 VBLKL8	53000	5.90	210000	7.70	232000	12.59

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8B
SEMIVOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): HH891228A08 Date Analyzed: 12/28/89
 Instrument ID: 08 Time Analyzed: 1217

	IS1 (DCB) AREA #	RT	IS2 (NPT) AREA #	RT	IS3 (ANT) AREA #	RT
12 HOUR STD	121000	6.78	370000	8.55	197000	11.10
UPPER LIMIT	242000		740000		394000	*
LOWER LIMIT	60500		185000		98500	
EPA SAMPLE NO.						
01 B201A	133000	6.77	336000	8.54	193000	11.09
02 SBLK44	163000	6.73	520000	8.49	259000	11.05

IS1 (DCB) = 1,4-Dichlorobenzene-d4
 IS2 (NPT) = Naphthalene-d8
 IS3 (ANT) = Acenaphthene-d10
 UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8C
SEMIVOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): HH891228A08 Date Analyzed: 12/28/89
 Instrument ID: 08 Time Analyzed: 1217

	IS4 (PHN) AREA #	RT	IS5 (CRY) AREA #	RT	IS6 (PRY) AREA #	RT
12 HOUR STD	250000	13.27	162000	17.22	128000	19.67
UPPER LIMIT	500000		324000		256000	
LOWER LIMIT	125000		81000		64000	
EPA SAMPLE NO.						
01 B201A	205000	13.29	165000	17.27	188000	19.77
02 SBLK44	320000	13.25	200000	17.22	161000	19.69

IS4 (PHN) = Phenanthrene-d10
 IS5 (CRY) = Chrysene-d12
 IS6 (PRY) = Perylene-d12

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8B
SEMIVOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): HH900110C08 Date Analyzed: 01/10/90
 Instrument ID: 08 Time Analyzed: 0746

	IS1 (DCB) AREA #	RT	IS2 (NPT) AREA #	RT	IS3 (ANT) AREA #	RT
=====	=====	=====	=====	=====	=====	=====
12 HOUR STD	168000	7.65	539000	9.32	289000	11.74
=====	=====	=====	=====	=====	=====	=====
UPPER LIMIT	336000		1078000		578000	
=====	=====	=====	=====	=====	=====	=====
LOWER LIMIT	84000		269500		144500	
=====	=====	=====	=====	=====	=====	=====
EPA SAMPLE NO.						
=====	=====	=====	=====	=====	=====	=====
01 B202TAR	168000	7.67	571000	9.32	307000	11.72
02 B202TARMS	186000	7.75	611000	9.40	318000	11.80
03 B202TARMSD	204000	7.75	670000	9.40	351000	11.77
04 SBLK79	207000	7.67	618000	9.32	331000	11.74

IS1 (DCB) = 1,4-Dichlorobenzene-d4
 IS2 (NPT) = Naphthalene-d8
 IS3 (ANT) = Acenaphthene-d10

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8C
SEMIVOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): HH900110C08 Date Analyzed: 01/10/90
 Instrument ID: 08 Time Analyzed: 0746

	IS4 (PHN) AREA #	RT	IS5 (CRY) AREA #	RT	IS6 (PRY) AREA #	RT
12 HOUR STD	409000	13.75	302000	17.50	266000	20.74
UPPER LIMIT	818000		604000		532000	
LOWER LIMIT	204500		151000		133000	
EPA SAMPLE NO.						
01 B202TAR	458000	13.75	306000	17.49	291000	20.72
02 B202TARMS	451000	13.82	316000	17.60	290000	20.90
03 B202TARMSD	494000	13.80	310000	17.55	273000	20.84
04 SBLK79	451000	13.77	282000	17.52	215000	20.77

IS4 (PHN) = Phenanthrene-d10
 IS5 (CRY) = Chrysene-d12
 IS6 (PRY) = Perylene-d12

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8B
SEMIVOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): HG891221B20 Date Analyzed: 12/21/89
 Instrument ID: 20 Time Analyzed: 2305

	IS1 (DCB) AREA #	RT	IS2 (NPT) AREA #	RT	IS3 (ANT) AREA #	RT
=====	=====	=====	=====	=====	=====	=====
12 HOUR STD	258000	6.33	725000	7.98	349000	10.37
=====	=====	=====	=====	=====	=====	=====
UPPER LIMIT	516000		1450000		698000	
=====	=====	=====	=====	=====	=====	=====
LOWER LIMIT	129000		362500		174500	
=====	=====	=====	=====	=====	=====	=====
EPA SAMPLE NO.						
=====	=====	=====	=====	=====	=====	=====
01 B201ADL	269000	6.33	858000	7.97	349000	10.37
02 B202A	229000	6.35	667000	7.98	284000	10.35
03 B202B	213000	6.33	754000	7.98	396000	10.37
04 B202BDL	227000	6.32	677000	7.95	355000	10.32
05 B202C	306000	6.35	837000	7.98	421000	10.37
06 B201ADLMS	274000	6.27	909000	7.93	426000	10.32
07 B201ADLMSD	238000	6.32	837000	7.97	401000	10.35
08 SBLK44	344000	6.27	884000	7.92	394000	10.30

IS1 (DCB) = 1,4-Dichlorobenzene-d4 UPPER LIMIT = + 100%
 IS2 (NPT) = Naphthalene-d8 of internal standard area.
 IS3 (ANT) = Acenaphthene-d10 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8C
SEMIVOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): HG891221B20 Date Analyzed: 12/21/89
 Instrument ID: 20 Time Analyzed: 2305

	IS4 (PHN) AREA #	RT	IS5 (CRY) AREA #	RT	IS6 (PRY) AREA #	RT
=====	=====	=====	=====	=====	=====	=====
12 HOUR STD	386000	12.37	227000	16.00	177000	18.57
=====	=====	=====	=====	=====	=====	=====
UPPER LIMIT	772000		454000		354000	
=====	=====	=====	=====	=====	=====	=====
LOWER LIMIT	193000		113500		88500	
=====	=====	=====	=====	=====	=====	=====
EPA SAMPLE NO.						
=====	=====	=====	=====	=====	=====	=====
01 B201ADL	402000	12.37	187000	15.99	118000	18.54
02 B202A	326000	12.35	158000	15.95	100000	18.45
03 B202B	488000	12.37	315000	15.95	248000	18.47
04 B202BDL	417000	12.30	234000	15.87	205000	18.37
05 B202C	442000	12.37	264000	15.95	213000	18.47
06 B201ADLMS	493000	12.32	291000	15.92	223000	18.42
07 B201ADLMSD	473000	12.35	336000	15.95	245000	18.45
08 SBLK44	399000	12.30	197000	15.95	127000	18.49

IS4 (PHN) = Phenanthrene-d10
 IS5 (CRY) = Chrysene-d12
 IS6 (PRY) = Perylene-d12

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

8B
SEMIVOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): HG900102C21 Date Analyzed: 01/02/90
 Instrument ID: 21 Time Analyzed: 0239

	IS1(DCB) AREA #	RT	IS2(NPT) AREA #	RT	IS3(ANT) AREA #	RT
12 HOUR STD	142000	7.50	446000	9.22	247000	11.69
UPPER LIMIT	284000		892000		494000	
LOWER LIMIT	71000		223000		123500	
EPA SAMPLE NO.						
01 B201B	243000	7.47	824000	9.19	433000	11.67
02 SBLK82	172000	7.48	480000	9.20	246000	11.67

IS1 (DCB) = 1,4-Dichlorobenzene-d4 UPPER LIMIT = + 100%
 IS2 (NPT) = Naphthalene-d8 of internal standard area.
 IS3 (ANT) = Acenaphthene-d10 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

SEMIVOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): HG900102C21 Date Analyzed: 01/02/90
 Instrument ID: 21 Time Analyzed: 0239

	IS4 (PHN) AREA #	RT	IS5 (CRY) AREA #	RT	IS6 (PRY) AREA #	RT
12 HOUR STD	390000	13.77	331000	17.65	257000	21.17
UPPER LIMIT	780000		662000		514000	
LOWER LIMIT	195000		165500		128500	
EPA SAMPLE NO.						
01 B201B	593000	13.75	386000	17.60	334000	21.09
02 SBLK82	362000	13.75	309000	17.60	252000	21.07

IS4 (PHN) = Phenanthrene-d10
 IS5 (CRY) = Chrysene-d12
 IS6 (PRY) = Perylene-d12

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

II. SAMPLE DATA PACKAGE

The Sample Data Package shall include data for analyses of all samples in one Sample Delivery Group, including field samples, reanalyses, blanks, matrix spikes and matrix spike duplicates. The Sample Data Package consists of the following:

1. CASE NARRATIVE
2. TRAFFIC REPORTS
3. VOLATILES DATA
4. SEMIVOLATILES DATA
5. PESTICIDE / PCB DATA

CASE#: 18756 SDG#: 07 SAS#: _____

1. CASE NARRATIVE

This document shall be clearly labeled " Case Narrative " and shall contain : laboratory name ; sample numbers in the Sample Delivery Group (SDG), differentiating between initial analyses and re - analyses ; SDG number ; Contract number ; and detailed documentation of any quality control, sample, shipment and / or analytical problems encountered in processing the samples reported in the data package.

Whenever data from sample re - analyses are submitted, the Contractor shall state in the Case Narrative for each re - analysis, whether it considers the re - analysis to be billable, and if so, why.

The contractor must also include documentation of any internal quality control processes used, a summary of corrective actions taken, and the resolution.

CLP CASE NARRATIVE--CASE # 18756
Contract No. 2-88 (REVS) SDG No. 07
CompuChem Laboratories, Inc.

Sample Numbers: B201A, B201B, B202A, B202B, B202C, B202TAR

This portion of Case # 18756 consisted of six solid samples for volatile and semivolatile analysis. The samples were received intact on 12/20/80 via Federal Express in properly sealed shipping containers with traffic reports. Due to a discrepancy between the sample container and the chain-of-custody, the C-O-C ID was used for sample B202A. Moisture content of the samples ranged from 9% to 64%. The pH values were within the values specified in EPA protocols.

VOLATILES:

Because of the nature of the samples, it was not possible to analyze the Low Level sample spikes within holding time requirements. All other volatile fractions were extracted and analyzed within holding time requirements. Because of the level of sample constituents, two analyses, a Low Level and a Medium Level, were reported for samples B201A, B201B, and B202TAR. The Low Level analyses of those three samples all contained toluene, 2-hexanone, and total xylenes at levels outside their analytical ranges. In addition, sample B201A also contained benzene and ethylbenzene and sample B202TAR, methylene chloride, acetone, and ethylbenzene, at levels outside their analytical range in the Low Level analyses. Because of sample inhomogeneity and the differences in preparation procedures for the Low and Medium analyses, some of the reported concentrations did not compare well between the two reported analyses for these three samples. The total xylene concentration was mistakenly flagged with an "E" in the dilution of sample B202TAR. Since the analysis partially separates the xylenes, the individual xylene concentrations were acceptable. The surrogate DB-toluene failed recovery criteria in sample B201A due to matrix interference. All other surrogate recoveries met acceptance criteria. Both the Low and Medium Level matrix spike/matrix spike duplicate results were acceptable. Due to a scheduling error, the sample requested for use as the spike original was not used. Toluene failed recovery criteria in the Medium Level MS and MSD, and also failed %RPD in the Medium Level analyses.

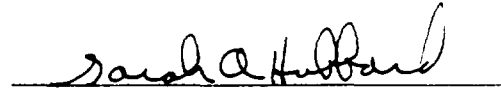
SEMIVOLATILES:

All semivolatile fractions were extracted and analyzed within holding time requirements. Coeluting compounds were flagged with an "X" on Form I. Because of the levels of 2-methylphenol, 4-methylphenol, and 2,4-dimethylphenol present, two analyses were reported for sample B201A, a 20:1 and a 50:1 dilution. Both a neat and a 5:1 analysis were reported for sample B202B because of the levels of 4-methylphenol and 2,4-dimethylphenol present. In addition to the above-mentioned TCL's, such compounds as phenol and various polyanomatic hydrocarbons were detected in one or more of these samples. Due to the viscosity of the extract, only a 5:1 dilution of the Medium Level extraction was reported for sample B201B. Because of the dilution necessary, no surrogate recovery data was available for the Low Level matrix spike/matrix spike duplicate. The recovery for DS-nitrobenzene in sample B201A exceeded recovery criteria. All other surrogates met recovery criteria. Because of the dilution necessary, there was no Low Level spike recovery data. A Low Level blank spike, extracted and analyzed with these samples, had acceptable results.

The failure of the %RPD for 2,4-dinitrotoluene in the blank spike was allowed since the %RPD criteria are internal CompuChem requirements and not contractual necessities. The Medium Level matrix spike/matrix spike duplicate results were acceptable. The recoveries for n-nitroso-di-n-propylamine and 1,2,4-trichlorobenzene failed QC limits for both the MS and MSD. In addition, the recovery for 1,4-dichlorobenzene failed QC limits in the MSD. The %RPD exceeded QC limits for 1,4-dichlorobenzene, 1,2,4-trichlorobenzene, and acenaphthene.

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or his designee as verified by the following signature.

Note: This report was paginated for reference and accountability in decreasing numerical sequence.



Sarah A. Hubbard 1/18/90
Technical Reviewer



DETECTION LIMIT CALCULATION CLARIFICATION

To protect our GC columns from unnecessary contamination, soil samples prepared according to Caucus Protocol methods are routinely diluted 5:1. Through a series of experiments we have determined that our instrument Detection Limit for pesticides is 5X lower than the EPA Contract Required Quantitation Limit (CRQL). We, therefore, only adjust our detection limits if the dilution necessary to analyze the sample is greater than 5:1. If the sample is diluted by a factor of X the detection limit is adjusted by X/5 instead of X.

Doug McCormack

Doug McCormack
Manager, GC Laboratory

Robert E. Meierer

Robert E. Meierer
Vice - President, Quality Assurance



**CHANGES TO THE EPA'S ORGANIC STATEMENT OF
WORK (SOW) FOR THE CONTRACT LABORATORY PROGRAM (CLP)**

Effective with samples received at CompuChem Laboratories, Inc. on April 3, 1989, the new "SOW for Organics Analysis; Multi-Media, Multi-Concentration" will be in effect. The new SOW is dated 2/88 with revisions dated 9/88.

EPA introduces modifications to the CLP SOW for a variety of reasons. They are:

- as a result of technical caucuses attended by representatives from EPA regional laboratories and the CLP, new or modified analytical methods are required,
- as a result of analytical data being supplied to the Agency by the CLP laboratories, QC acceptance criteria are updated and made a requirement of the program,
- as a result of requirements by the end users of the data (the EPA regions and the Program Office), changes to the deliverable requirements of the CLP are necessary.

As a service to our clients utilizing the EPA CLP SOW for their analytical needs, the following information is provided to point out the substantive changes between the new SOW and the previous one (10/86 with revisions through 8/87).

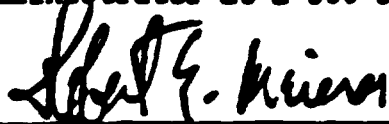
"KEY" CHANGES TO THE ORGANIC SOW

- 1) Wide bore capillary columns (internal diameter greater than 0.32 mm) are allowed for pesticide/PCB analysis in addition to packed columns. Fused silica capillary columns remain an optional confirmation column.

- 2) The Sample Data Summary Package (SDSP) was introduced in the 10/86 SOW. The SDSP consists of copies of specific summary forms. In the 2/88 SOW, Form VIII, the Internal Standard Area Summary Form for volatile and semi-volatile analyses, has been added to the SDSP. Form VIII for volatiles and semi-volatiles is also to be included in the QC Summary Package.
- 3) A clarification for dealing with the three xylene isomers has been added. For total xylenes, where three isomers are quantified as two peaks, the calibration range of each peak should be considered separately, e.g., a diluted analysis is not required for total xylenes unless the concentration of either peak exceeds 200 ug/l.
- 4) The volatile target analyte, 2-butanone, is to be quantified against the first internal standard, bromochloromethane. In the prior SOW, 2-butanone was quantified against the second internal standard, 1,4-difluorobenzene.
- 5) Turnaround time for the delivery of data has been lowered from 40 days to 35 days for routine analytical service (RAS) work associated with the Superfund program. The turnaround time is calculated from the Validated Time of Sample Receipt (VTSR) of the last sample in the Sample Delivery Group (SDG).

The above represents the major changes in the new organic SOW for EPA's CLP. There are, of course, other minor changes which have not been addressed in this announcement. Of note, the hardcopy deliverable forms have not changed and, in fact, retain the same revision dates as the prior SOW.

If there are any questions concerning the information presented, please feel free to contact your account administrator at 1-800-833-8097.



Robert E. Meierer
Vice President of Quality Assurance

DATA REPORTING QUALIFIERS

For reporting results to EPA, the following result qualifiers are used. Additional flags or footnotes explaining results are encouraged. However, the definition of each flag must be explicit.

VALUE - If the result is a value greater than or equal to the detection limit, report the value.

U - Indicates compound was analyzed but not detected. The sample quantitation limit must be corrected for dilution and for percent moisture. For example, 10 U for phenol in water if the sample final volume is the protocol-specified final volume. If a 1 to 10 dilution of extract is necessary, the reported limit is 100 U. For a soil sample, the value must also be adjusted for percent moisture. For example, if the sample had 24% moisture and a 1 to 10 dilution factor, the sample quantitation limit for phenol (330 U) would be corrected to:

$$\frac{(330 \text{ U}) \times df}{D} \quad \text{where } D = \frac{100 - \% \text{ moisture}}{100}$$

and df = dilution factor

$$\text{at } 24\% \text{ moisture, } D = \frac{100 - 24}{100} = 0.76$$

$$\frac{(330 \text{ U}) \times 10}{.76} = 4300 \text{ U rounded to the appropriate number of significant figures}$$

For soil sample subjected to GPC clean-up procedures, the ERL is also multiplied by 2, to account for the fact that only half of the extract is recovered.

J - Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed, or when the mass spectral data indicate the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than zero. For example, if the sample quantitation limit is 10 ug/L, but a concentration is 3 ug/L is calculated, report it as 3J. The sample quantitation limit must be adjusted for both dilution and percent moisture as discussed for the U flag, so that if a sample with 24% moisture and a 1 to 10 dilution factor has a calculated concentration of 300 ug/L and a sample quantitation limit of 430 ug/kg, report the concentration as 300J on Form 1.

C - This flag applies to pesticide results where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ug/ml in the final extract shall be confirmed by GC/MS.

DATA REPORTING QUALIFIERS - PAGE 2

- B - This flag is used when the analyte is found in the associated blank as well as in the sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action. This flag must be used for a TIC as well as for a positively identified TCL compound.**

- E - This flag identifies compounds whose concentrations exceed the calibration range of the GC/MS instrument for that specific analysis. This flag will not apply to pesticides/PCBs analyzed by GC/EC methods. If one or more compounds have a response greater than full scale, the sample or extract must be diluted and re-analyzed according to the specifications. All such compounds with a response greater than full scale should have the concentration flagged with an "E" on the Form I for the original analysis. If the dilution of the extract causes any compounds identified in the first analysis to be below the calibration range in the second analysis, then the results of both analyses shall be reported on separate Forms I. The Form I for the diluted sample shall have the "DL" suffix appended to the sample number.**

- D - This flag identifies all compounds identified in an analysis at a secondary dilution factor. If a sample or extract is re-analyzed at a higher dilution factor, as in the "E" flag above, the "DL" suffix is appended to the sample number on the Form I for the diluted sample and all concentration values reported on that Form I are flagged with the "D" flag.**

- A - This flag indicates that TIC is a suspected aldol-condensation product.**

- X - Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the Sample Data Summary Package and the Case Narrative. If more than one is required, use "Y" and "Z", as needed. If more than five qualifiers are required for a sample result, use the "X" flag to combine several flags, as needed. For instance, the "X" flag might combine the "A", "B", and "D" flags for some sample.**



COMPUCHEM LABORATORIES

QUALITY ASSURANCE NOTICE

With the advent of the new organics Statement of Work (SOW 2/88, Revision: 9/88) participants in EPA's Contract Laboratory Program (CLP) are required to provide hard copy and diskette deliverables. CompuChem employs the Finnigan QA Formaster Program (Format A) to generate these requirements using data files from our analytical instrumentation. Currently, and independently, quantitation reports are generated by the instruments and are used with CompuChem-developed software to calculate results. The GC and GC/MS quantitation routines employ the convention of carrying at least one extra significant figure until the mathematical computations are completed. Then, the quantitative results are rounded to the SOW-required number of significant figures for reporting. In addition, the algorithm used by the Formaster Program is slightly different than that employed in CompuChem's software routines. Therefore, results presented in the supportive data supplied with our deliverables packages may be slightly different than those which appear on the hard copy forms generated via Formaster.

This notice serves to alert the end users of these data packages as to the reason why slight differences may be observed.

Robert E. Meierer
Director of Quality Assurance

QUALITY ASSURANCE NOTICE

Specific guidelines are presented in the EPA CLP Organic Statement of Work for the positive qualitative identification of compounds through mass spectral interpretation. Applying these guidelines absolutely may not be possible when the nature of the sample is less than pure reference material. Where the mass spectral pattern of a compound to be identified demonstrates interferences or resolution from one or more additional compounds, either unknowns, internal standards, or surrogate standards, the "+" sign is added to the top of the dual spectra page.

Linda Fowler 4/10/89


Linda Fowler
Sr. Quality Assurance Specialist

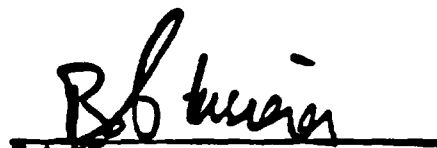
Robert E. Malerer 4/12/89

Robert E. Malerer
Vice President, Quality Assurance

LABORATORY NOTICE

On June 15, 1985 CompuChem Laboratories began adding D3-2,4-Dinitrophenol to all standards and samples. The purpose of this addition is to enable the Laboratory to have higher and more consistent analytical sensitivity for the native 2,4-Dinitrophenol. The peak corresponding to the deuterated analog is clearly labeled on each RIC as D3/1 and will not be searched and reported as a tentatively identified compound (TIC). This compound is not being used as an internal or surrogate standard.


L. Richard Flynn,
Development Chemist


Bob Meierer,
Director of Quality Assurance

2. TRAFFIC REPORTS

A copy of the Sample Traffic Reports in Item A for all of the samples in the SDG. The Traffic Reports shall be arranged in increasing EPA number order, considering both letters and numbering in ordering samples.

3. VOLATILES DATA

A. QC SUMMARY

B. SAMPLE DATA

C. STANDARDS DATA

D. RAW QC DATA

CASE#: 18756 SDG#: 07 SAS#: _____

A. QC SUMMARY

- (1) **Surrogate Percent Recovery Summary (Form II VOA)**
- (2) **Matrix Spike / Matrix Spike Duplicate Summary (Form III VOA)**
- (3) **Method Blank Summary (Form IV VOA)**

**(If more than a single form is necessary , forms must
be arranged in chronological order by date of analysis
of the blank)**

- (4) **GC / MS Tuning and Mass Calibration (Form V VOA)**

BFB in chronological order ; by instrument

- (5) **Internal Standard Area Summary (Form VIII VOA)**

In chronological order ; by instrument

(1) Surrogate Percent Recovery Summary (Form II VOA)

2B
SOIL VOLATILE SURROGATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Level: (low/med) LOW

	EPA SAMPLE NO.	S1 (TOL) #	S2 (BFB) #	S3 (DCE) #	OTHER	TOT OUT
01	B201A	80 *	90	97		1
02	B201B	90	99	94		0
03	B202A	95	92	97		0
04	B202B	92	96	97		0
05	B202C	94	91	97		0
06	B202TAR	88	93	97		0
07	B202AMS	101	98	92		0
08	B202AMSD	114	113	102		0
09	VBLKL8	107	98	95		0
10	VBLKB4	97	87	97		0
11	VBLKF5	92	91	92		0

QC LIMITS

S1 (TOL) = Toluene-d8 (81-117)
 S2 (BFB) = Bromofluorobenzene (74-121)
 S3 (DCE) = 1,2-Dichloroethane-d4 (70-121)

Column to be used to flag recovery values
 * Values outside of contract required QC limits
 D Surrogates diluted out

2B
SOIL VOLATILE SURROGATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Level: (low/med) MED

	EPA SAMPLE NO.	S1 (TOL) #	S2 (BFB) #	S3 (DCE) #	OTHER	TOT OUT
01	B201ARE	91	93	111		0
02	B201BRE	115	105	113		0
03	B202TARDL	82	110	109		0
04	B201BREMS	90	100	90		0
05	B201BREMSD	86	101	104		0
06	VBLKH1	110	103	117		0
07	VBLKH9	101	104	117		0
08	VBLKI6	112	106	121		0

QC LIMITS

S1 (TOL) = Toluene-d8 (81-117)
 S2 (BFB) = Bromofluorobenzene (74-121)
 S3 (DCE) = 1,2-Dichloroethane-d4 (70-121)

Column to be used to flag recovery values

* Values outside of contract required QC limits

D Surrogates diluted out

(2) Matrix Spike / Matrix Spike Duplicate Summary (Form III VOA)

3B
SOIL VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix Spike - EPA Sample No.: B202A Level: (low/med) LOW

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
1,1-Dichloroethene	56.8	0	58.1	102	59-172
Trichloroethene	56.8	0	47.3	83	62-137
Benzene	56.8	0	55.2	97	66-142
Toluene	56.8	1.45	53.9	92	59-139
Chlorobenzene	56.8	0	52.0	92	60-133

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
1,1-Dichloroethene	56.8	53.5	94	8	22	59-172
Trichloroethene	56.8	47.5	84	-1	24	62-137
Benzene	56.8	55.1	97	0	21	66-142
Toluene	56.8	54.0	93	-1	21	59-139
Chlorobenzene	56.8	50.9	90	2	21	60-133

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 0 out of 5 outside limits

Spike Recovery: 0 out of 10 outside limits

COMMENTS: CLP ,1875,67, , ,SOIL,309688,VOLATILE, ,
TUNE: 0019 122289 0013

SOIL VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix Spike - EPA Sample No.: B201BRE Level: (low/med) MED

COMPOUND	SPIKE ADDED (ug/Kg)	SAMPLE CONCENTRATION (ug/Kg)	MS CONCENTRATION (ug/Kg)	MS % REC #	QC LIMITS REC.
1,1-Dichloroethene	7100	0	4730	67	59-172
Trichloroethene	7100	0	6950	98	62-137
Benzene	7100	2160	9740	107	66-142
Toluene	7100	15900	18900	42 *	59-139
Chlorobenzene	7100	0	6970	98	60-133

COMPOUND	SPIKE ADDED (ug/Kg)	MSD CONCENTRATION (ug/Kg)	MSD % REC #	% RPD #	QC LIMITS RPD	REC.
1,1-Dichloroethene	7100	4330	61	9	22	59-172
Trichloroethene	7100	6970	98	0	24	62-137
Benzene	7100	10500	118	-10	21	66-142
Toluene	7100	18000	30 *	33 *	21	59-139
Chlorobenzene	7100	7230	102	-4	21	60-133

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 1 out of 5 outside limits

Spike Recovery: 2 out of 10 outside limits

COMMENTS: CLP ,1875,67, , ,SOIL,309686,VOLATILE, ,
 TUNE: 0013 122989 1535

8A
VOLATILE INTERNAL STANDARD AREA SUMMARY

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Lab File ID (Standard): GS900102A19 Date Analyzed: 01/02/90
 Instrument ID: 19 Time Analyzed: 0942
 Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) CAP

	IS1 (BCM) AREA #	RT	IS2 (DFB) AREA #	RT	IS3 (CBZ) AREA #	RT
=====	=====	=====	=====	=====	=====	=====
12 HOUR STD	52100	5.90	198000	7.68	214000	12.54
=====	=====	=====	=====	=====	=====	=====
UPPER LIMIT	104200		396000		428000	
=====	=====	=====	=====	=====	=====	=====
LOWER LIMIT	26050		99000		107000	
=====	=====	=====	=====	=====	=====	=====
EPA SAMPLE NO.						
=====	=====	=====	=====	=====	=====	=====
01 B202AMS	53700	5.88	212000	7.72	234000	12.60
02 B202AMSD	59400	5.87	223000	7.68	245000	12.60
03 VBLKL8	53000	5.90	210000	7.70	232000	12.59

IS1 (BCM) = Bromochloromethane
 IS2 (DFB) = 1,4-Difluorobenzene
 IS3 (CBZ) = Chlorobenzene

UPPER LIMIT = + 100%
 of internal standard area.
 LOWER LIMIT = - 50%
 of internal standard area.

Column used to flag internal standard area values with an asterisk

B. SAMPLE DATA

Sample data shall be arranged in packets with the Organic Analysis Data Sheet (Form I VOA , including Form I VOA - TIC), followed by the raw data for volatile samples. These sample packets should then be placed in increasing EPA number order, considering both letters and numbers in ordering samples.

- . **TCL Results - Organic Analysis Data Sheet (Form I VOA).**
Tabulated results (identification and quantitation) of the specified target compounds (Exhibit C).
- . **Tentatively Identified Compounds (Form I VOA - TIC).**
This form must be included even if no compounds are found. If so, indicate this on the form by entering " 0 " in the field for " Number Found ".
- . **Reconstructed total ion chromatograms (RIC) for each sample,**
sample extract, standard, blank and spiked sample.
- . **For each sample, by each compound identified.**
 - (a) **Copies of raw spectra and copies of background - subtracted mass spectra of target compounds listed in Exhibit C (TCL) that are identified in the sample and corresponding background - subtracted TCL standard mass spectra. Compound names must be clearly marked on all spectra.**
 - (b) **Copies of mass spectra of nonsurrogate organic compounds not listed in Exhibit C (TCL) (Tentatively Identified Compounds) which associated best - match spectra (three best matches).**

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009679C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 Date Analyzed: 12/22/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	(ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	11	U
74-83-9	-----Bromomethane	11	U
75-01-4	-----Vinyl Chloride	11	U
75-00-3	-----Chloroethane	11	U
75-09-2	-----Methylene Chloride	32	B
67-64-1	-----Acetone	150	B
75-15-0	-----Carbon Disulfide	5	U
75-35-4	-----1,1-Dichloroethene	5	U
75-34-3	-----1,1-Dichloroethane	5	U
540-59-0	-----1,2-Dichloroethene (total)	5	U
67-66-3	-----Chloroform	5	U
107-06-2	-----1,2-Dichloroethane	5	U
78-93-3	-----2-Butanone	120	U
71-55-6	-----1,1,1-Trichloroethane	5	U
56-23-5	-----Carbon Tetrachloride	5	U
108-05-4	-----Vinyl Acetate	11	U
75-27-4	-----Bromodichloromethane	5	U
78-87-5	-----1,2-Dichloropropane	5	U
10061-01-5	-----cis-1,3-Dichloropropene	5	U
79-01-6	-----Trichloroethene	5	U
124-48-1	-----Dibromochloromethane	5	U
79-00-5	-----1,1,2-Trichloroethane	5	U
71-43-2	-----Benzene	260	E
10061-02-6	-----Trans-1,3-Dichloropropene	5	U
75-25-2	-----Bromoform	5	U
108-10-1	-----4-Methyl-2-Pentanone	17	U
591-78-6	-----2-Hexanone	580	E
127-18-4	-----Tetrachloroethene	5	U
79-34-5	-----1,1,2,2-Tetrachloroethane	5	U
108-88-3	-----Toluene	720	E
108-90-7	-----Chlorobenzene	5	U
100-41-4	-----Ethylbenzene	420	E
100-42-5	-----Styrene	210	E
1330-20-7	-----Total Xylenes	1300	E

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201A

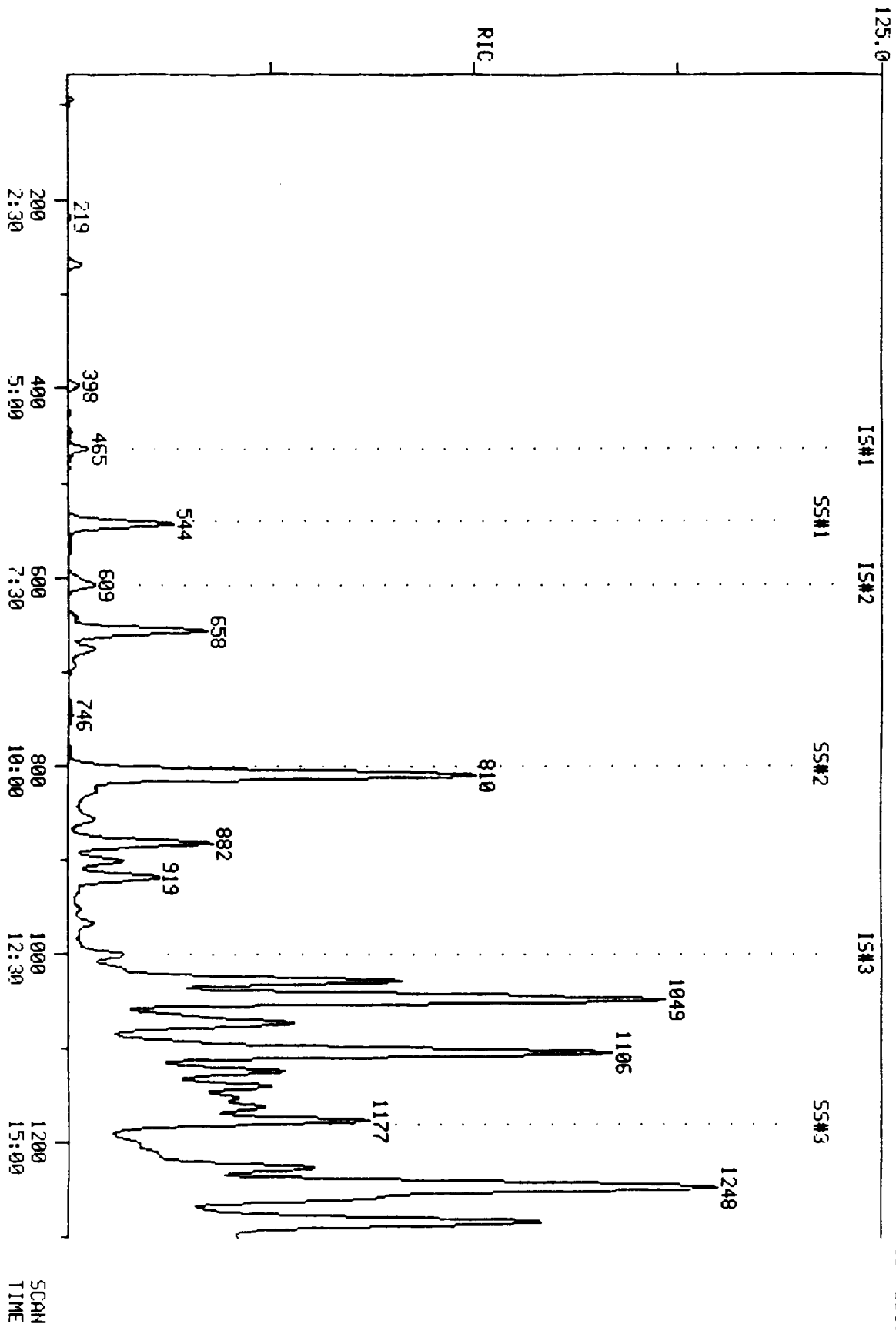
Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009679C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 9 Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 625-86-5	FURAN, 2,5-DIMETHYL-	8.22	240	J
2. 1192-62-7	ETHANONE, 1-(2-FURANYL)-	11.02	250	J
3.	UNKNOWN	13.44	460	J
4.	UNKNOWN	13.45	220	J
5.	UNKNOWN	14.05	190	J
6.	UNKNOWN	14.55	300	J
7. 100-66-3	BENZENE, METHOXY-	14.70	510	J
8. 103-65-1	BENZENE, PROPYL-	15.34	190	J
9.	UNKNOWN	15.60	1500	J
10.	SUBSTITUTED BENZENE	16.05	800	J

COMPUchem LABS
COMPUchem DATA: GH09679C19 SCANS 69 TO 1300
12/22/89 4:18:00
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
COND.S.:

1247990.



QUANTITATION REPORT FILE: GH009679C19
 DATA: GH009679C19.TI
 12/22/89 4:18:00
 SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A DN#19
 CONDS: :
 SUBMITTED BY: 19 ANALYST: 1422

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
 RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1
2	221 CHLOROMETHANE <74-87-3> RO#2
3	231 VINYL CHLORIDE <75-01-4> RO#3
4	220 BROMOMETHANE <78-83-9> RO#4
5	209 CHLOROETHANE <75-00-3> RO#5
6	216 1,1-DICHLOROETHENE <75-35-4> RO#8
7	254 CARBON DISULFIDE <75-15-0> RO#9
8	252 ACETONE (2-PROPANONE) <67-64-1> RO#13
9	*249 1,4-DIFLUOROBENZENE (IS) <540-36-3> RO#14
10	222 METHYLENE CHLORIDE <75-09-2> RO#16
11	226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17
12	214 1,1-DICHLOROETHANE <75-34-3> RO#19
13	257 VINYL ACETATE <108-05-4> RO#20
14	237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21
15	253 2-BUTANONE <78-93-3> RO#22
16	211 CHLOROFORM <67-66-2> RO#23
17	227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24
18	206 CARBON TETRACHLORIDE <56-23-5> RO#25
19	203 BENZENE <71-43-2> RO#26
20	215 1,2-DICHLOROETHANE <107-06-2> RO#27
21	*270 D5-CHLOROBENZENE (IS) RO#29
22	229 TRICHLOROETHENE <79-01-6> RO#30
23	217 1,2-DICHLOROPROPANE <78-87-5> RO#31
24	212 BROMODICHLOROMETHANE <75-27-4> RO#33
25	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35
26	256 4-METHYL-2-PENTANONE <108-01-1> RO#36
27	225 TOLUENE <108-88-3> RO#37
28	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38
29	228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39
30	224 TETRACHLOROETHENE <127-18-4> RO#41
31	255 2-HEXANONE <591-78-6> RO#42
32	208 DIBROMOCHLOROMETHANE <124-48-1> RO#43
33	207 CHLOROBENZENE <108-90-7> RO#45
34	219 ETHYLBENZENE <100-41-4> RO#47
35	330 M,P-XYLENE <133-02-7> RO#48
36	239 O-XYLENE <133-02-7> RO#49
37	251 STYRENE <100-42-5> RO#50
38	205 BROMOFORM <75-25-2> RO#51
39	223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54
40	*258 D4-1,2-DICHLOROETHANE RO#57
41	*247 BROMOFLUOROBENZENE <460-00-4> RO#58
42	*233 D8-TOLUENE RO#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	464	5:48	1	1.000	A BB	34851.	50.000 UG/KG	1.34
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	219	2:44	1	0.472	A BB	24465.	132.645 UG/KG	3.56
9	114	609	7:37	9	1.000	A BB	134420.	50.000 UG/KG	1.34
10	84	269	3:22	1	0.580	A BB	33717.	29.437 UG/KG	0.79
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	448	5:36	1	0.966	A BB	8006.	107.176 UG/KG	2.88
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	544	6:48	9	0.893	A BB	561099.	233.122 UG/KG	6.26
20	62	NOT FOUND							
21	117	1000	12:30	21	1.000	A BB	158516.	50.000 UG/KG	1.34
22	130	NOT FOUND							
23	63	NOT FOUND							
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	801	10:01	21	0.801	A BB	12617.	15.507 UG/KG	0.42
27	92	810	10:07	21	0.810	A BB	1524640.	651.540 UG/KG	17.50
28	75	NOT FOUND							
29	97	NOT FOUND							
30	164	NOT FOUND							
31	43	919	11:29	21	0.919	A VB	324090.	523.670 UG/KG	14.07
32	129	NOT FOUND							
33	112	NOT FOUND							
34	106	1029	12:52	21	1.029	A BV	615073.	382.836 UG/KG	10.28
35	106	1049	13:07	21	1.049	A VB	1458110.	652.791 UG/KG	17.54
36	106	1104	13:48	21	1.104	A BB	1024930.	519.806 UG/KG	13.96
37	104	1108	13:51	21	1.108	A BB	663203.	190.361 UG/KG	5.11
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	540	6:45	1	1.164	A BB	70282.	48.323 UG/KG	1.30
41	95	1181	14:46	21	1.181	A VB	122608.	45.209 UG/KG	1.21
42	98	801	10:01	21	0.801	A BB	132677.	39.945 UG/KG	1.07

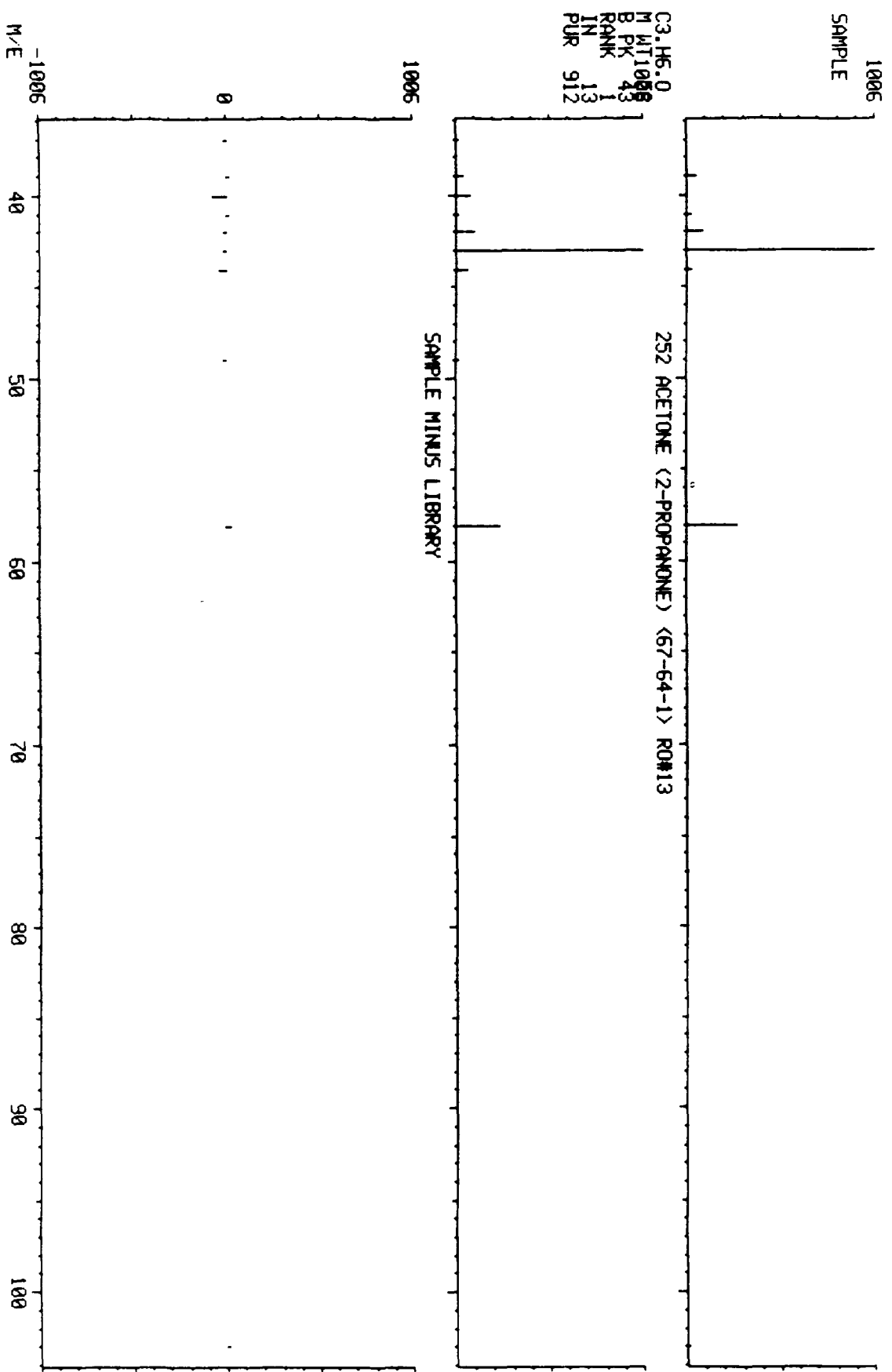
NO	RET(L)	RATID	RRT(L)	RATID	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:54	0.98	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:06		10.000			50.00		0.952	
3	1:11		10.000			50.00		1.256	
4	1:26		10.000			50.00		1.579	
5	1:34		10.000			50.00		0.796	
6	2:37		5.000			50.00		1.346	
7	2:46		5.000			50.00		3.709	
8	2:48	0.98	10.000	0.05	132.64	50.00	0.702	0.265	2.65
9	7:42	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	3:27	0.97	5.000	0.12	29.44	50.00	0.967	1.643	0.59
11	3:53		5.000			50.00		1.486	
12	4:34		5.000			50.00		2.635	
13	4:52		10.000			50.00		0.458	
14	5:33		5.000			50.00		1.675	

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:42	0.98	10.000	0.10	107.18	50.00	0.230	0.107	2.14
16	6:10		5.000			50.00		3.433	
17	6:18		5.000			50.00		0.814	
18	6:34		5.000			50.00		0.755	
19	6:53	0.99	5.000	0.18	233.12	50.00	4.174	0.895	4.66
20	6:58		5.000			50.00		2.523	
21	12:34	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	8:01		5.000			50.00		0.569	
23	8:22		5.000			50.00		0.352	
24	8:55		5.000			50.00		0.701	
25	9:41		5.000			50.00		0.690	
26	10:04	0.99	15.000	0.05	15.51	50.00	0.080	0.257	0.31
27	10:10	1.00	5.000	0.16	651.54	50.00	9.618	0.738	13.03
28	10:42		5.000			50.00		0.342	
29	10:58		5.000			50.00		0.349	
30	11:06		5.000			50.00		0.723	
31	11:33	0.99	15.000	0.06	523.67	50.00	2.045	0.195	10.47
32	11:35		5.000			50.00		0.690	
33	12:37		5.000			50.00		1.092	
34	12:55	1.00	5.000	0.21	382.84	50.00	3.880	0.507	7.66
35	13:10	1.00	5.000	0.21	652.79	50.00	9.199	0.705	13.06
36	13:52	1.00	5.000	0.22	519.81	50.00	6.466	0.622	10.40
37	13:54	1.00	5.000	0.22	190.36	50.00	4.184	1.099	3.81
38	14:07		5.000			50.00		0.664	
39	15:16		5.000			50.00		0.557	
40	6:52	0.98	5.000	0.23	48.32	50.00	2.017	2.087	0.97
41	14:49	1.00	5.000	0.24	45.21	50.00	0.773	0.855	0.90
42	10:04	0.99	5.000	0.16	39.95	50.00	0.837	1.048	0.80

LIBRARY SEARCH
12/22/89 4:18:00 + 2:44
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009679C19 # 219
BASE M/E: 43
RIC: 4463.



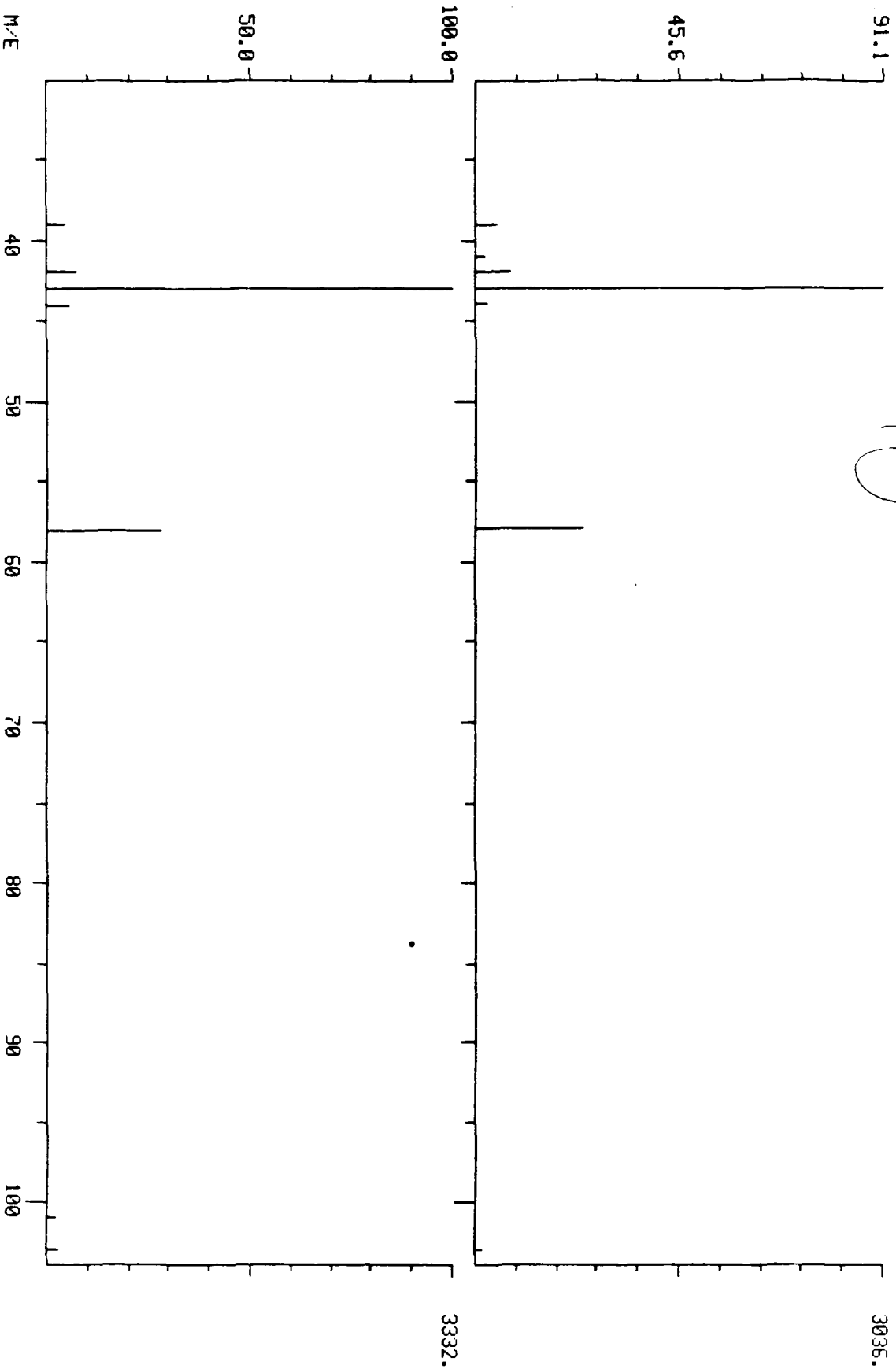
COMFUCHEM LABS

DATA: GH009679019 #219

BASE M/E: 43/ 43
RIC: 4463.7 5023.

1430

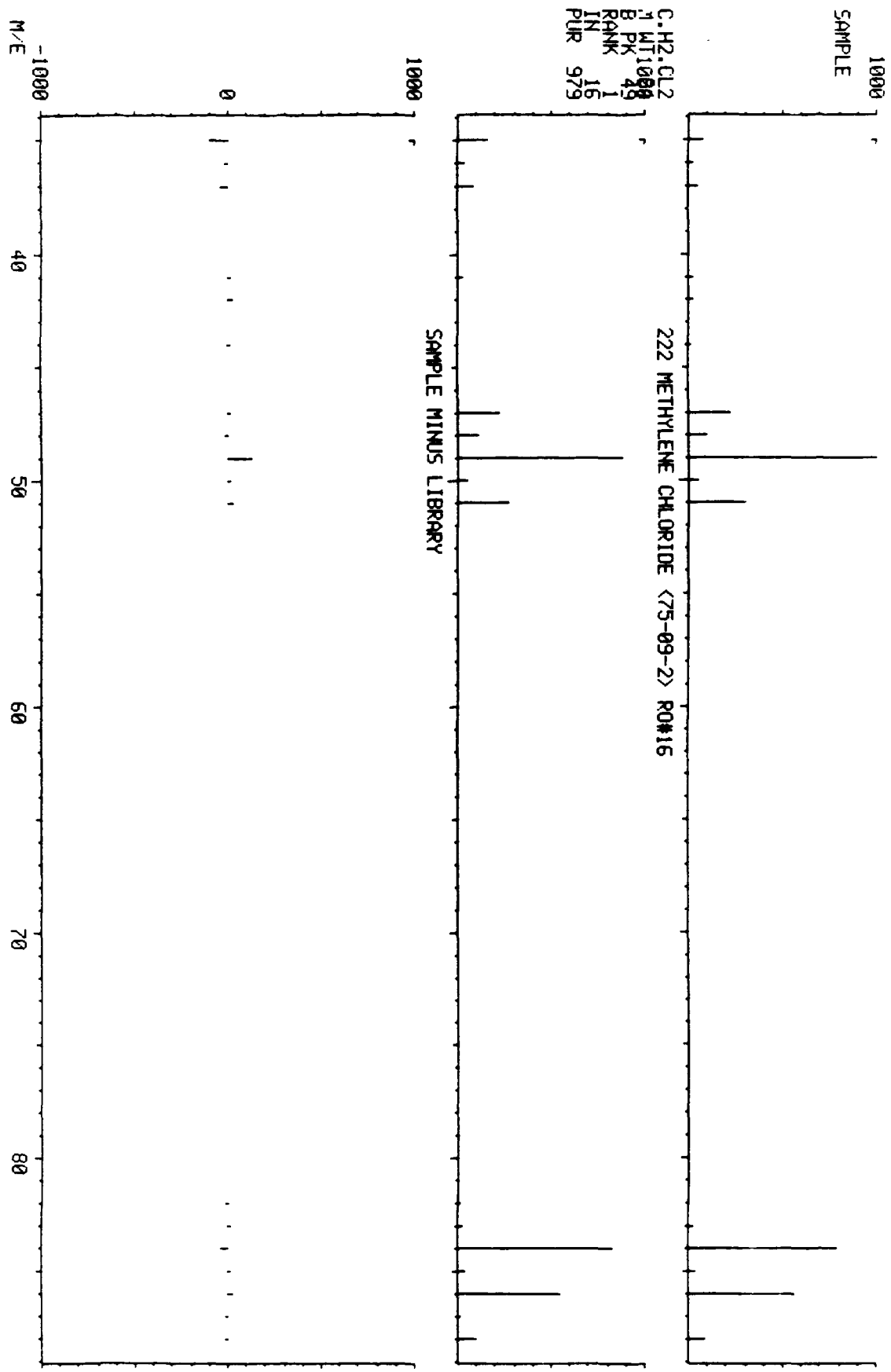
QUAL MASS SPECTRUM
12/22/89 4:18:00 + 2:44
SAMPLE: 5G CC#309679 CASE#18756.7 EPANB201A ON#19
ENHANCED (5 158 2N) (252) ACETONE (2-PROPANONE) (67-64-1) R0#13



COMPUchem LABS
LIBRARY SEARCH
12/22/89 4:18:00 + 3:22
SAMPLE: 5G OC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N 0T)

DATA: GH009679C19 # 269
BASE M/E: 49
RIC: 18079.

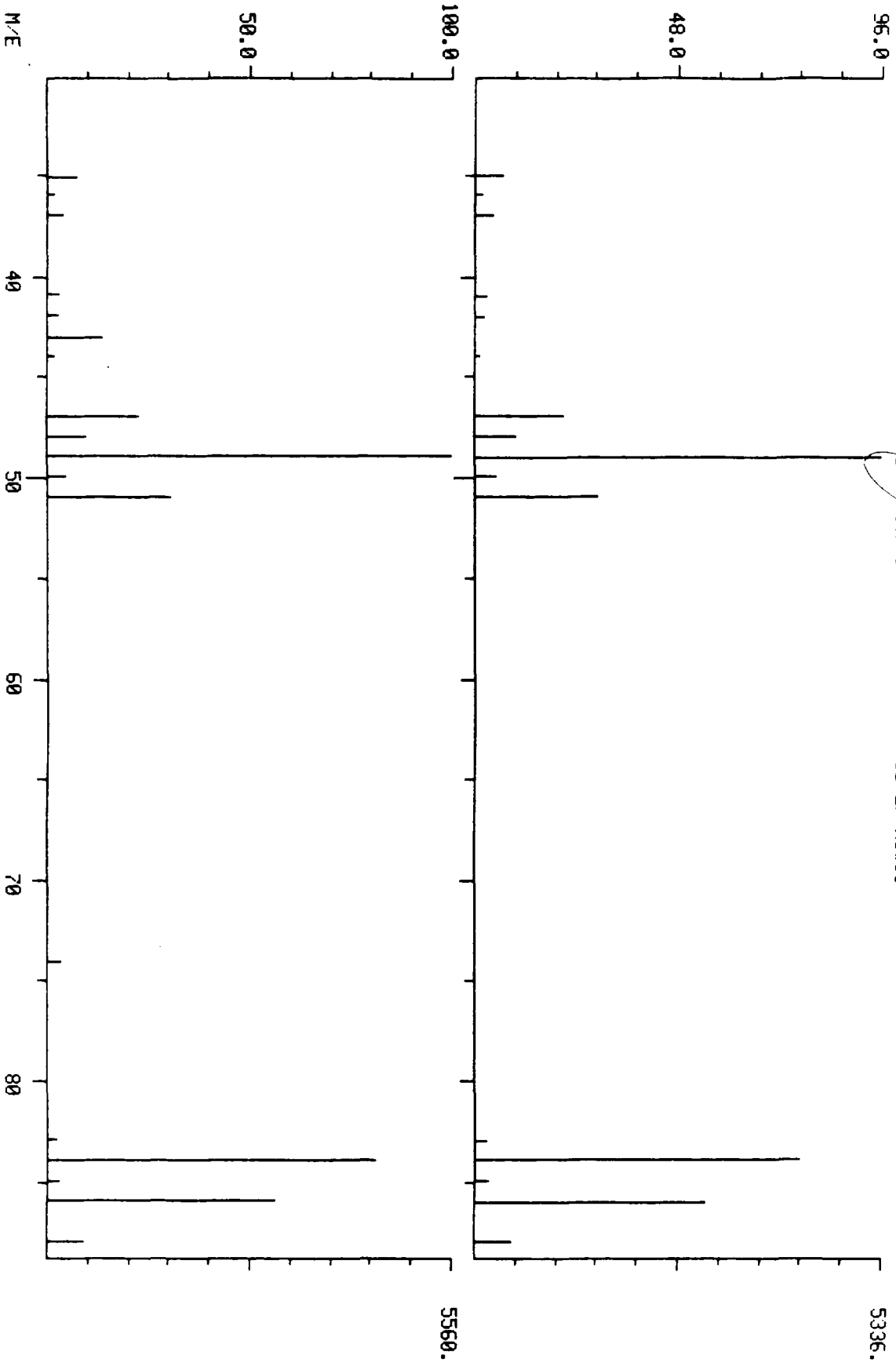
C.H2.CL2
1 MIT 1000
B PK 49
IN 1
IN 16
PUR 979



DUAL MASS SPECTRUM
12/22/89 4:18:00 + 3:22
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 15B 2N) 222 METHYLENE CHLORIDE <75-09-2> R0#16

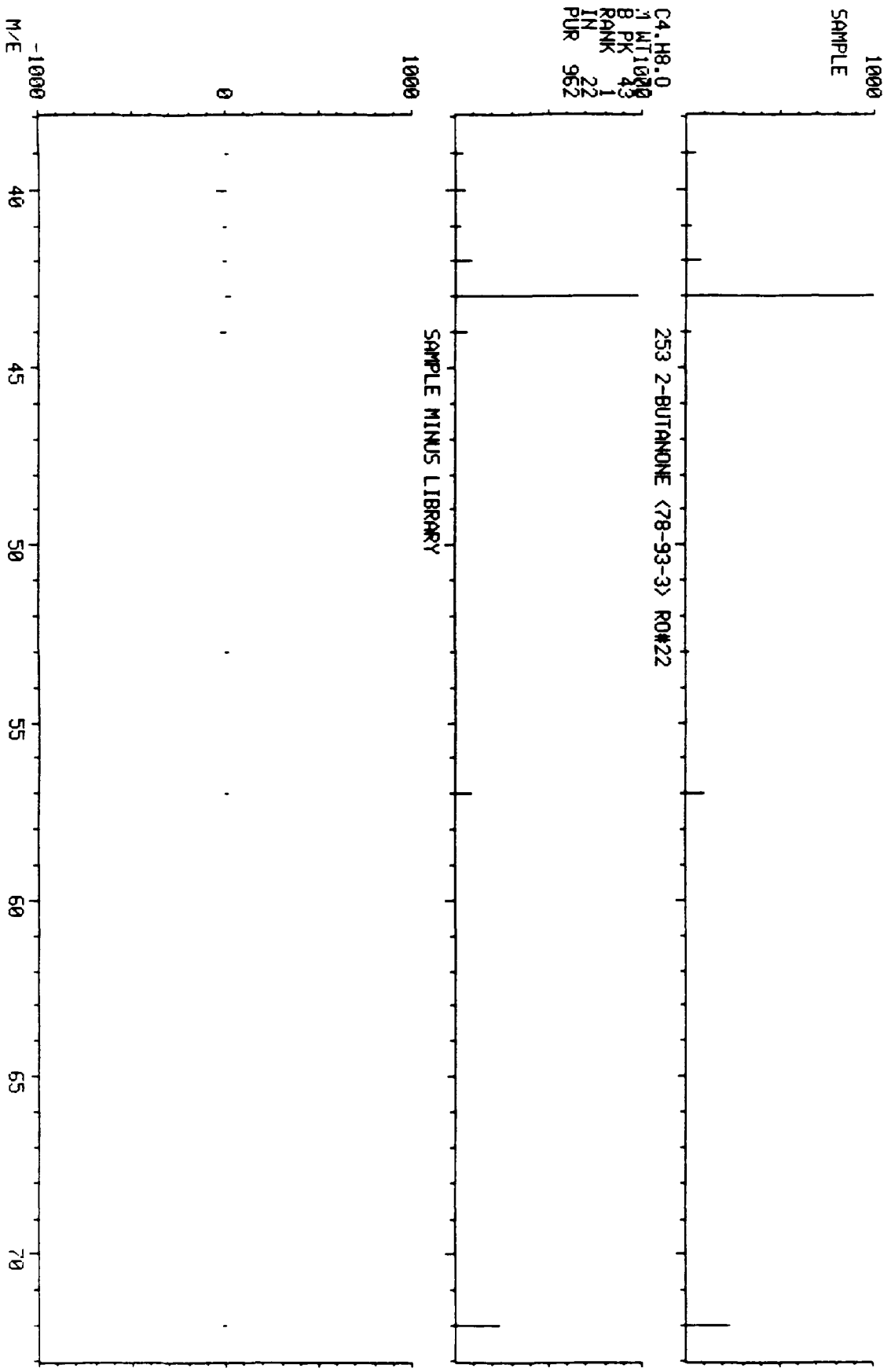
COMPUchem LABS

DATA: GH009679C19 #259 BASE M/E: 49/ 49
RIC: 18079. / 19807.



LIBRARY SEARCH
 12/22/89 4:18:00 + 5:36
 SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
 ENHANCED (5 15B 2N 0T)

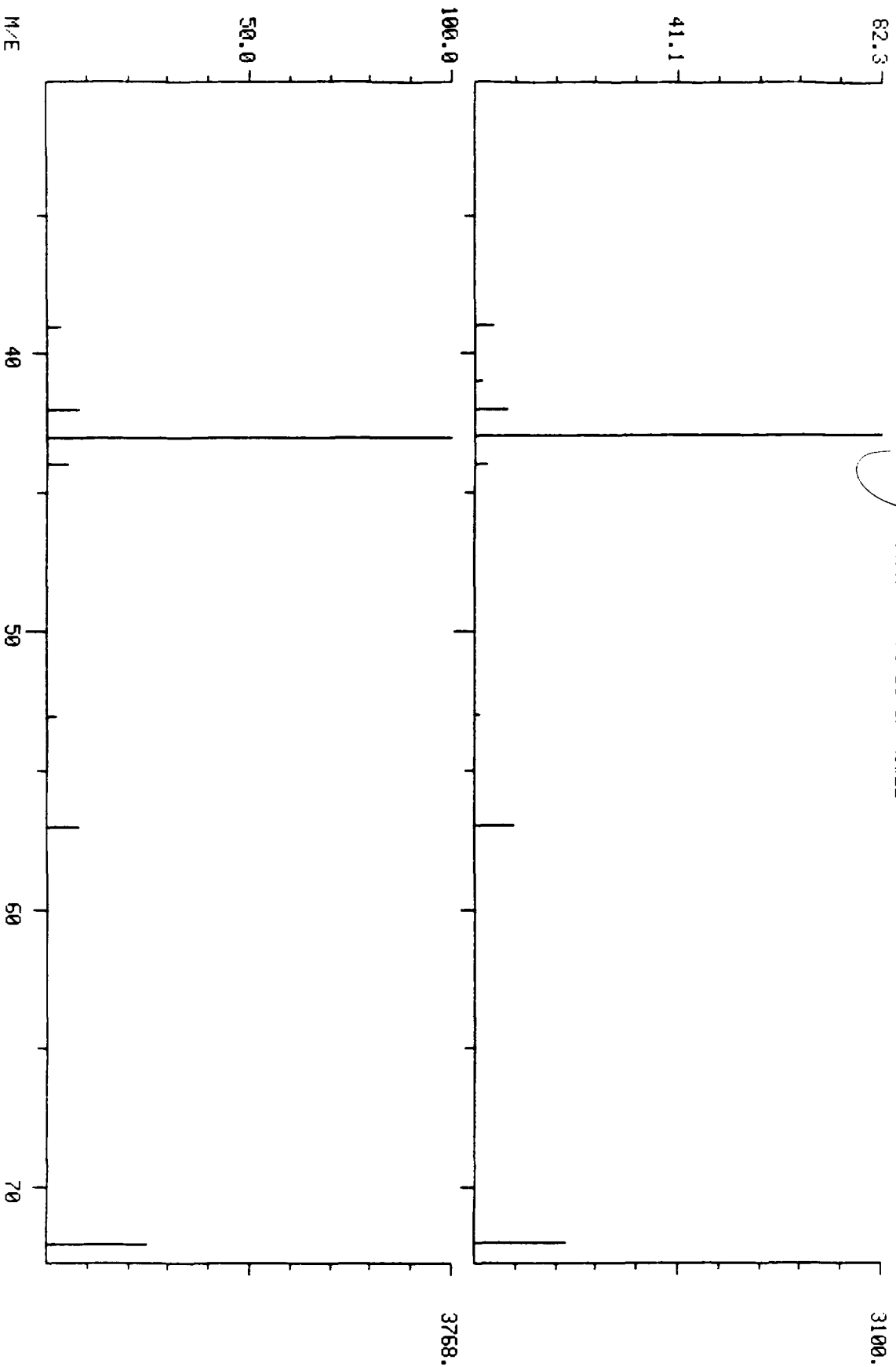
COMPUCHEN LABS
 DATA: GH009679C19 # 448
 BASE M/E: 43
 RIC: 4663.



DUAL MASS SPECTRUM
12/22/89 4:18:00 + 5:36
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2H) / 253 2-BUTANONE (78-93-3) RO#22

COMPUCHEM LABS

DATA: GH009679C19 #448 BASE M/E: 43/ 43
RIC: 4653./ 5703.



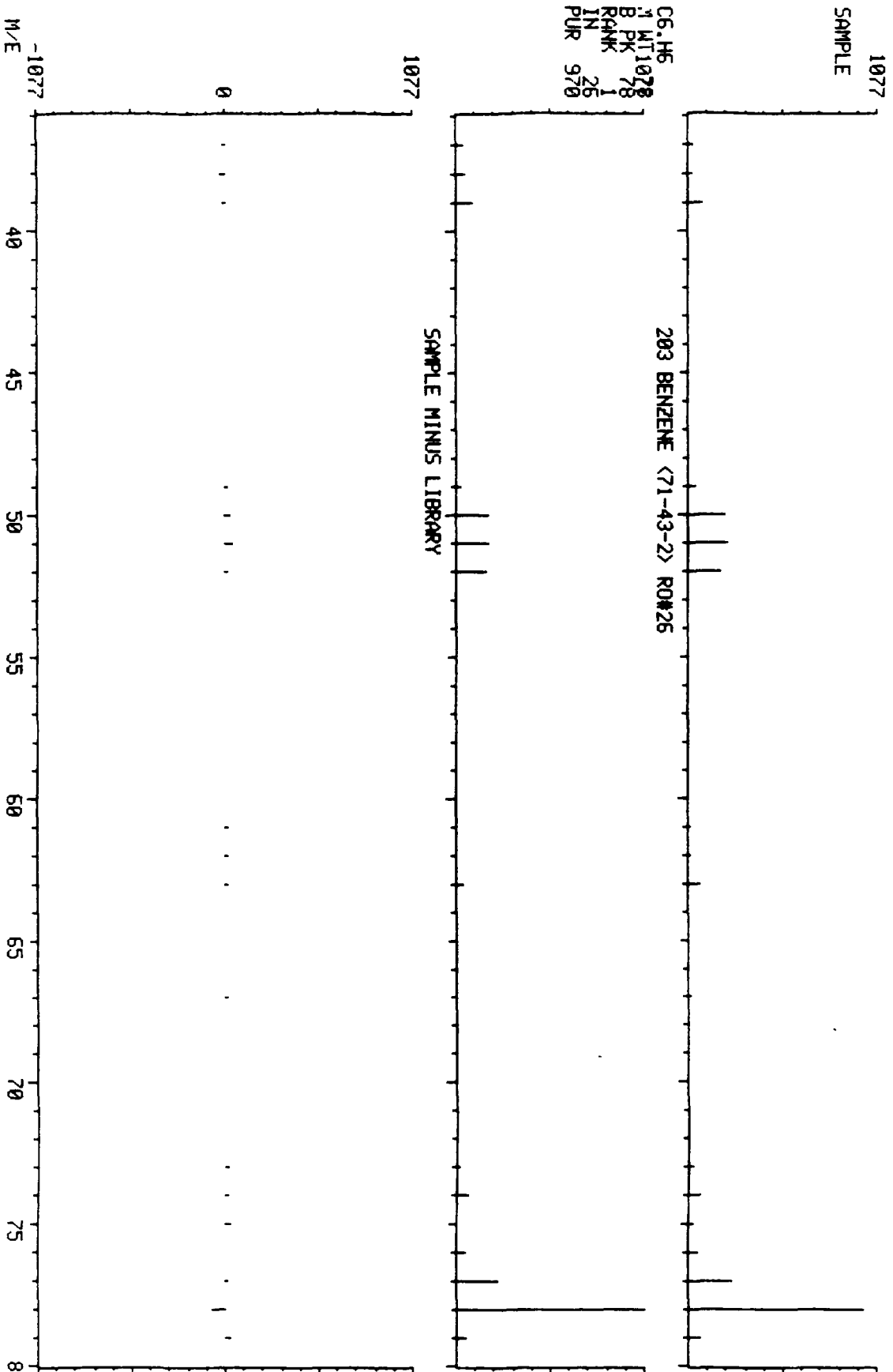
LIBRARY SEARCH
12/22/89 4:18:00 + 6:48
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A OH#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: CH009679C19 # 544

BASE M/E: 78
RIC: 123007.

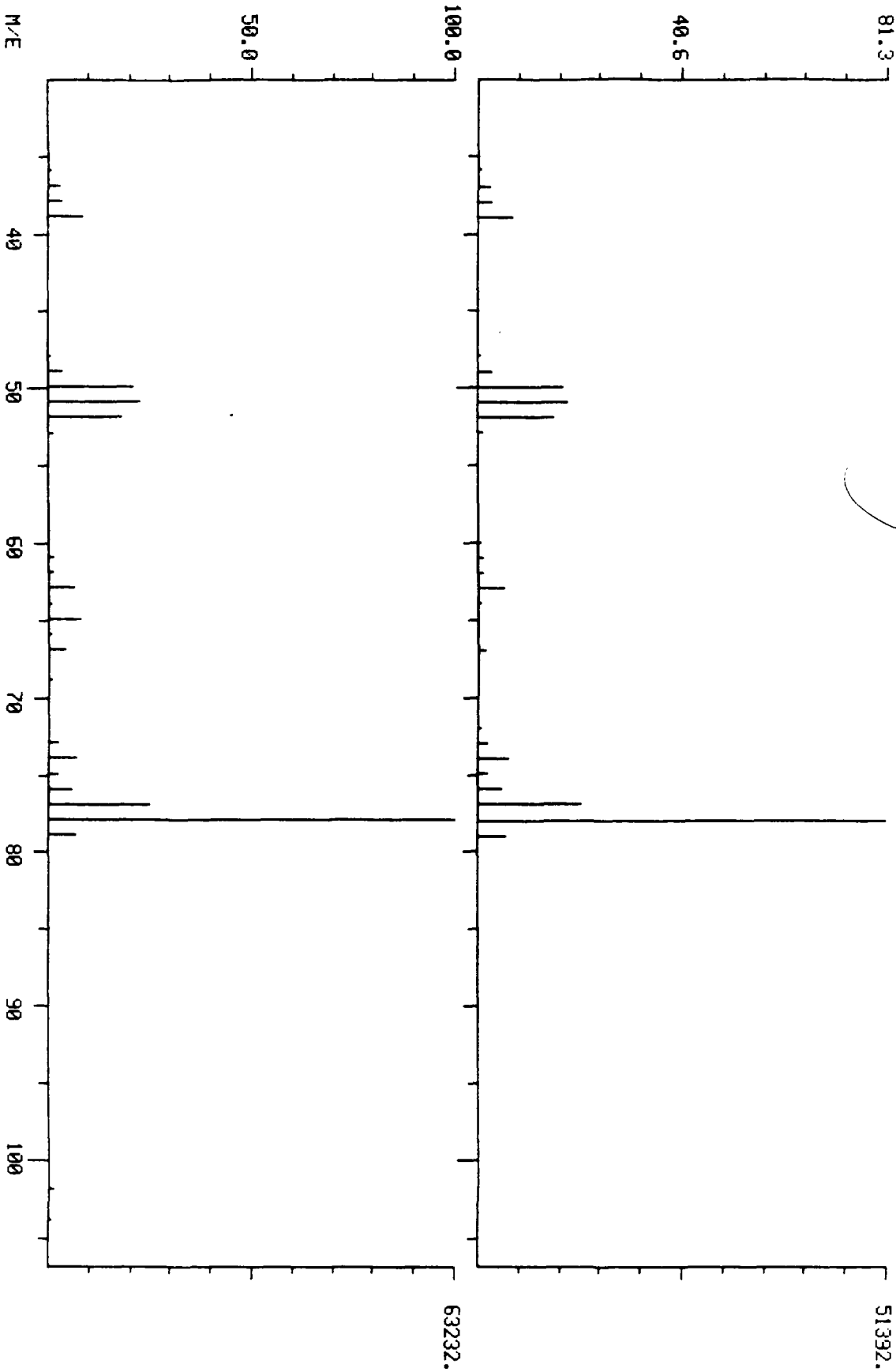
C6.H6
M WT 1078
B PK 78
RANK 1
IN 26
PUR 970



DUAL MASS SPECTRUM
12/22/89 4:18:00 + 6:48
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#8201A ON#19
ENHANCED (5 158 2N) 203 BENZENE (71-43-2) R0#26

CONFLUENT LABS

DATA: GH009679C19 #544 BASE M/E: 78/ 78
RIC: 124671. / 159999.



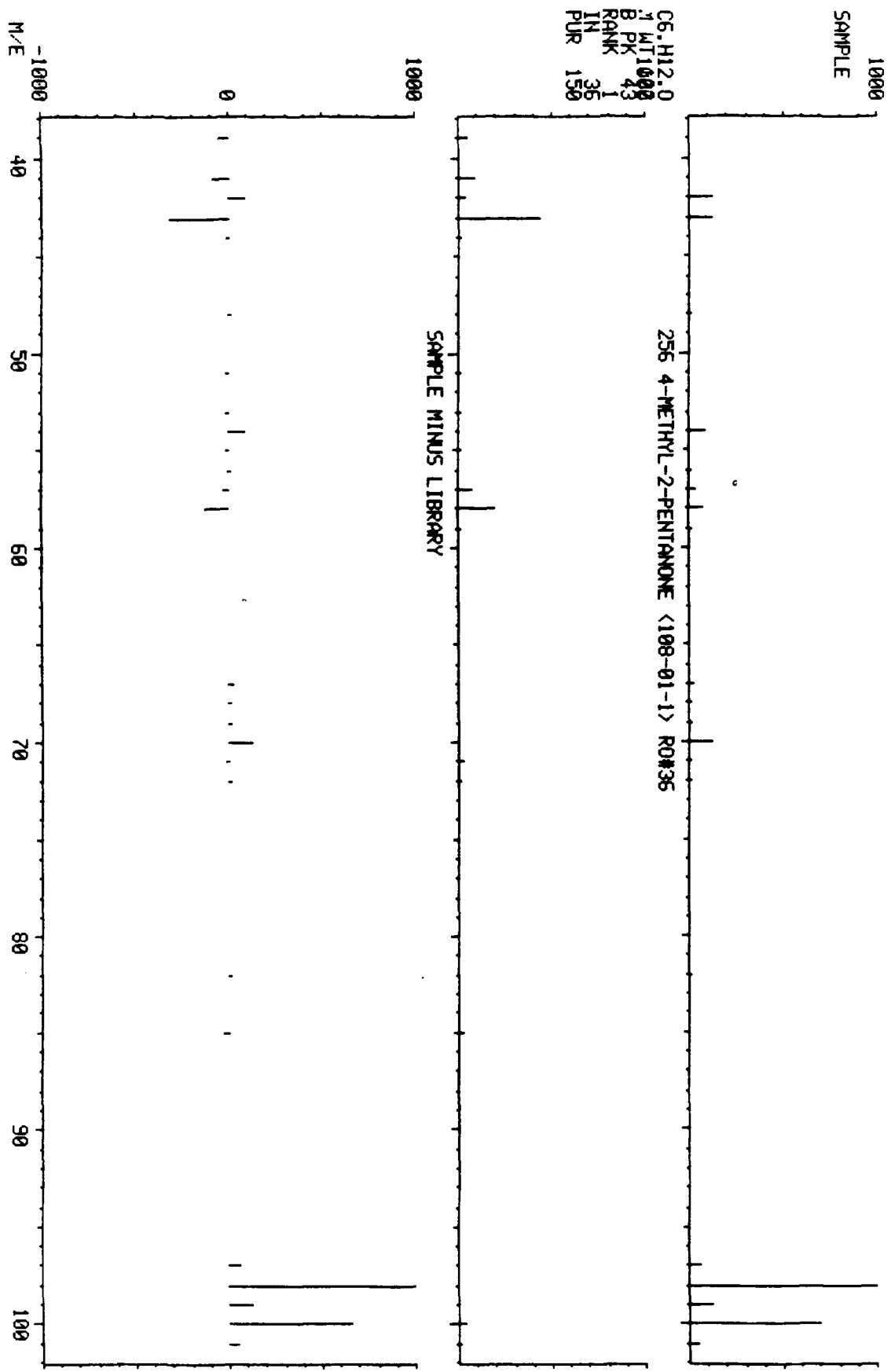


COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 4:18:00 + 10:01
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N 0T)

DATA: GH009679C19 # 801

BASE M/E: 98
RIC: 30463.

CG.H12.0
I.M.I.1000
B.PK 43
RANK 1
IN 35
PUR 150

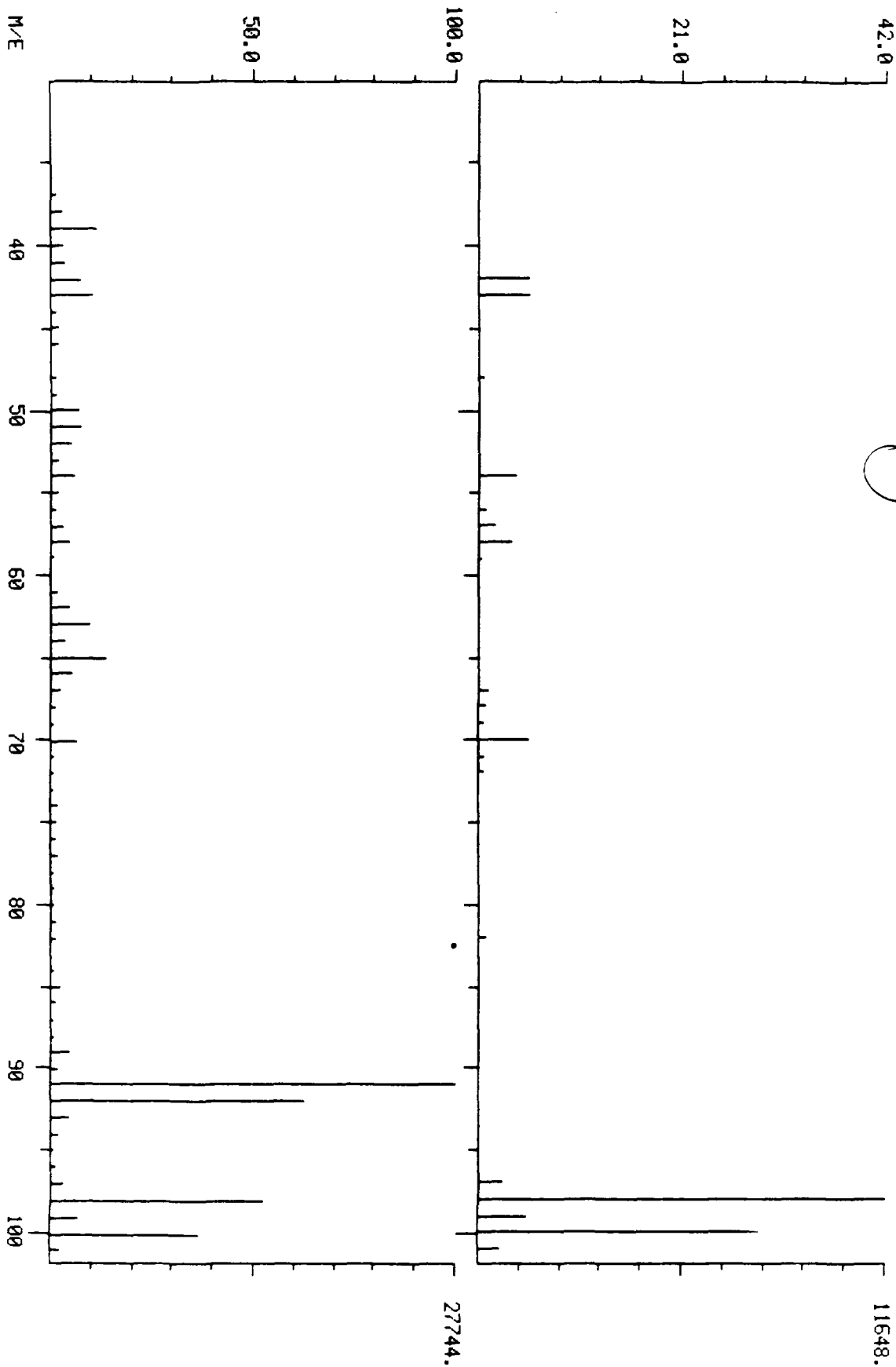


DUAL MASS SPECTRUM
12/22/89 4:18:00 + 10:01
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (S 158 2N) / 256 4-METHYL-2-PENTANONE <108-01-1> R0#36

COMPUCHEM LABS

DATA: GH09679C19 #801

BASE M/E: 98/ 91
RIC: 30463. / 117119.



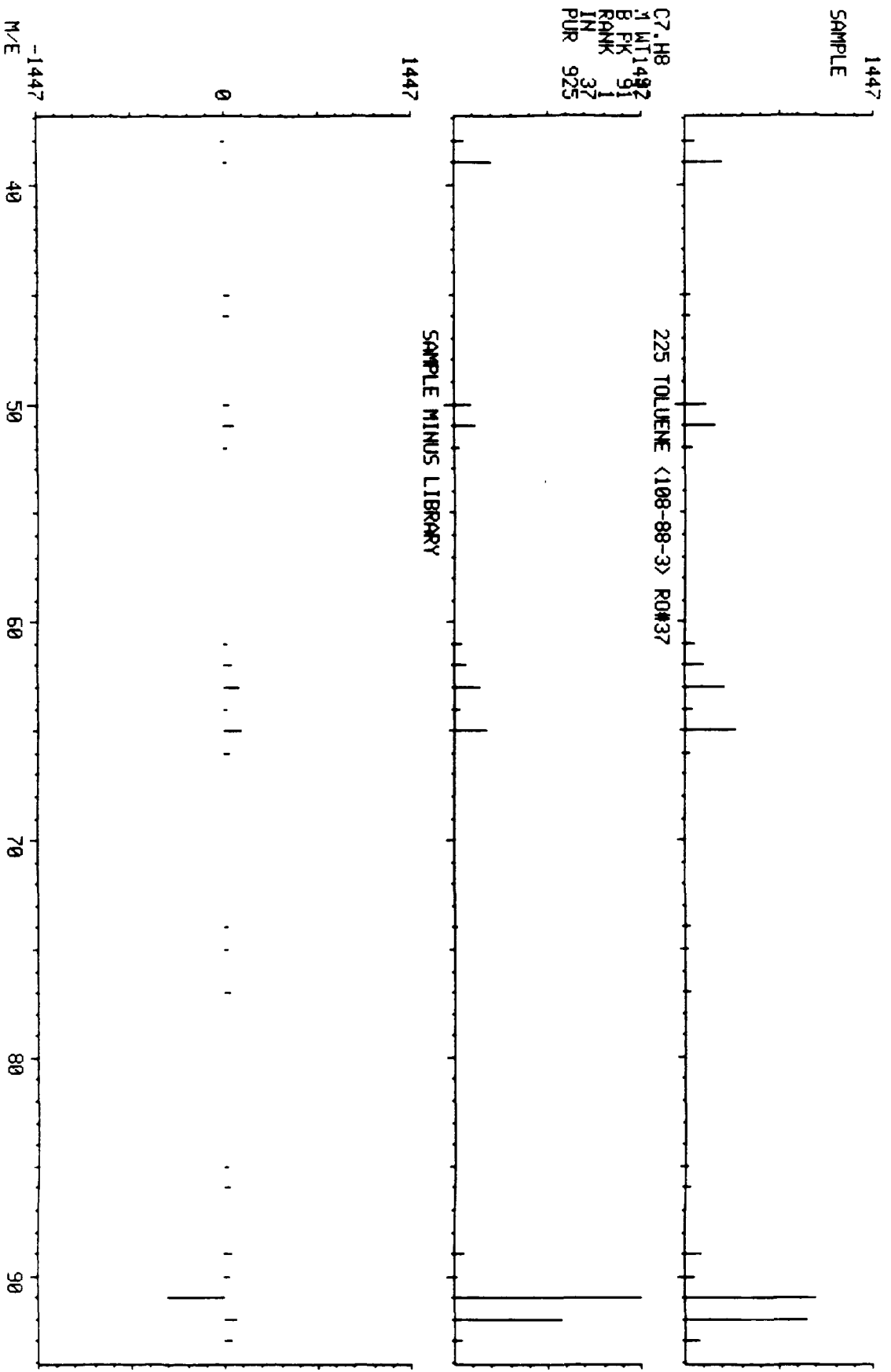
LIBRARY SEARCH
12/22/89 4:18:00 + 10:07
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N 0T)

COMPUchem LABS

DATA: GH009679C19 # 810

BASE M/E: 91
RIC: 432127.

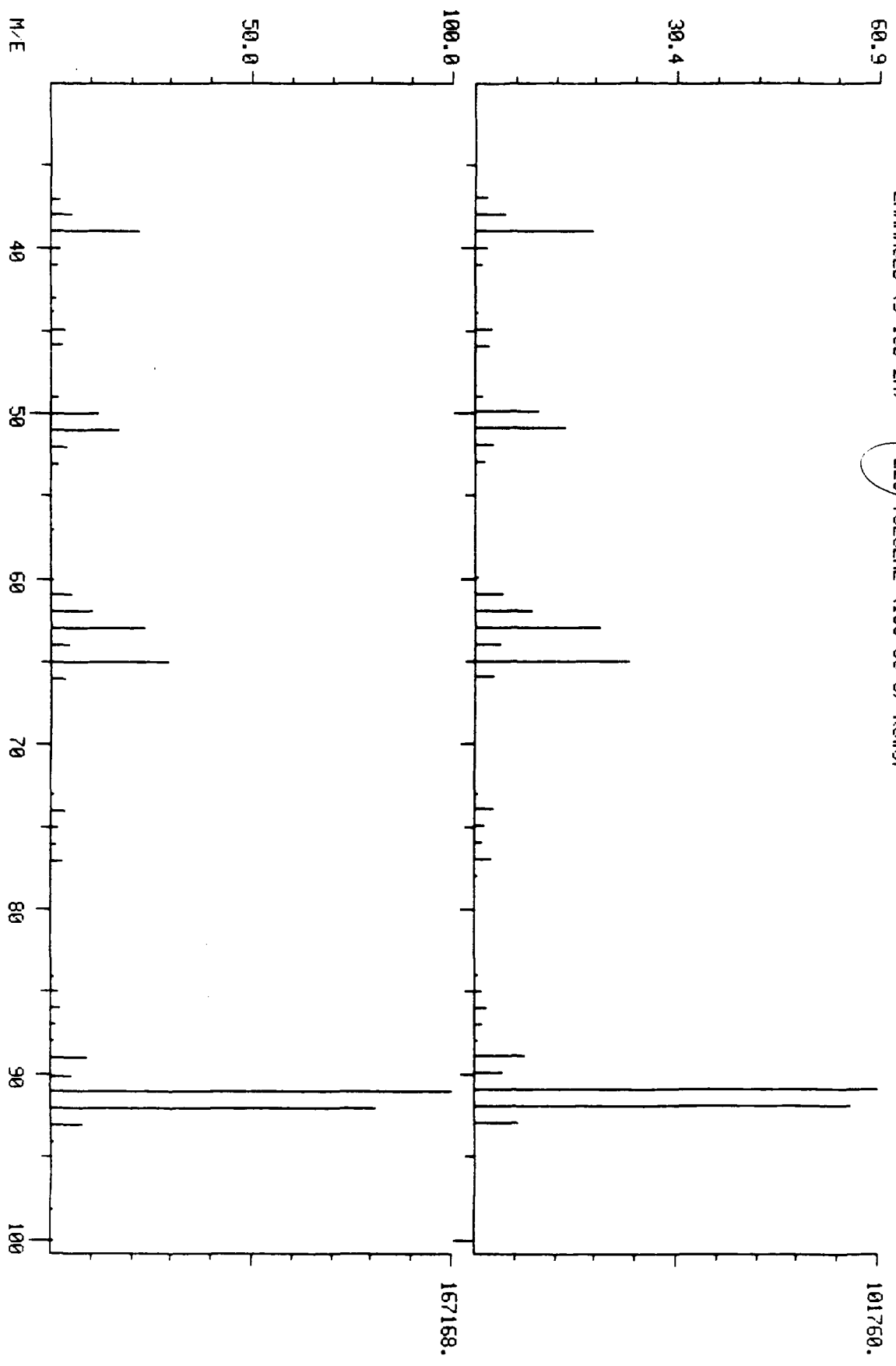
C7.H8
Y WT 1492
B PK 91
RANK 37
IN 1
PUR 925



DUAL MASS SPECTRUM
12/22/89 4:18:00 + 10:07
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2M) 225 TOLUENE <108-88-3> RM#37

COMPUCHEM LABS

DATA: GH009679C19 #810 BASE M/E: 91 / 91
RIC: 451071. / 625663.

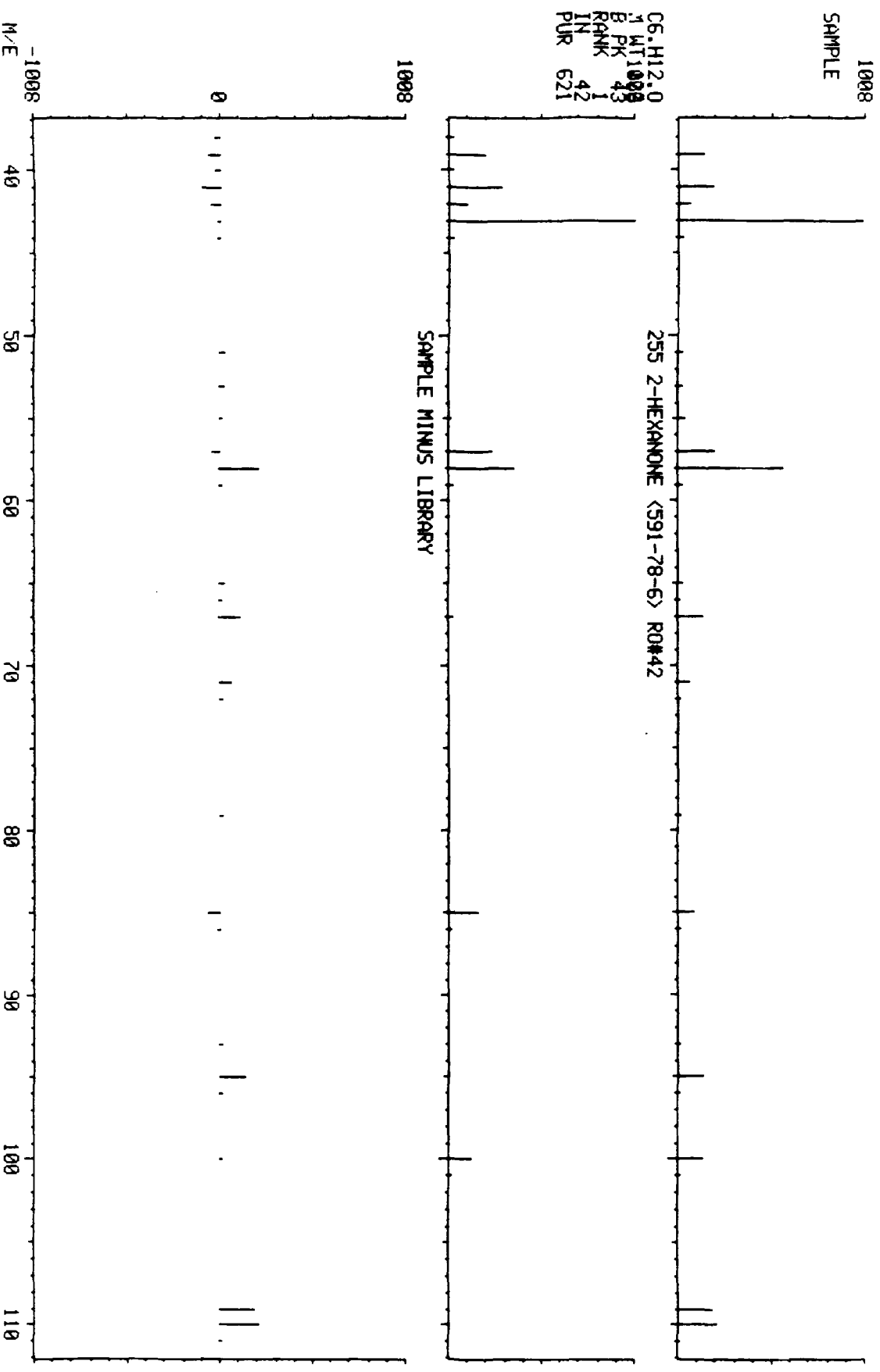


LIBRARY SEARCH
12/22/89 4:18:00 + 11:29
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 15B 2N 0T)

COMPUCHEN LABS

DATA: CH009679C19 # 919

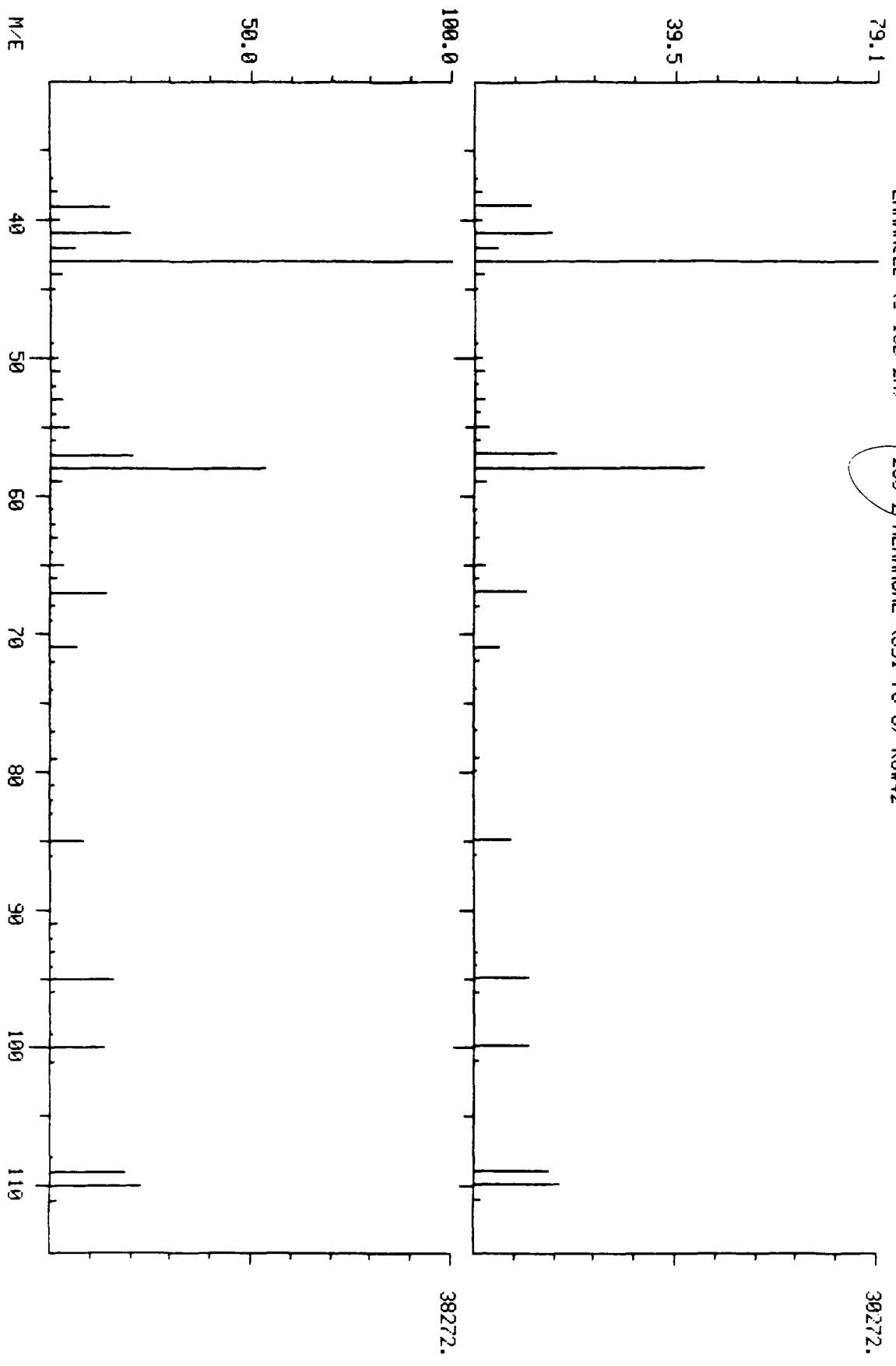
BASE M/E: 43
RIC: 101503.



DUAL MASS SPECTRUM
12/22/89 4:18:00 + 11:29
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#8201A DN#19
ENHANCED (5 158 2N) 255 2-HEXANONE (591-78-6) RQ#42

COMPUCHEM LABS

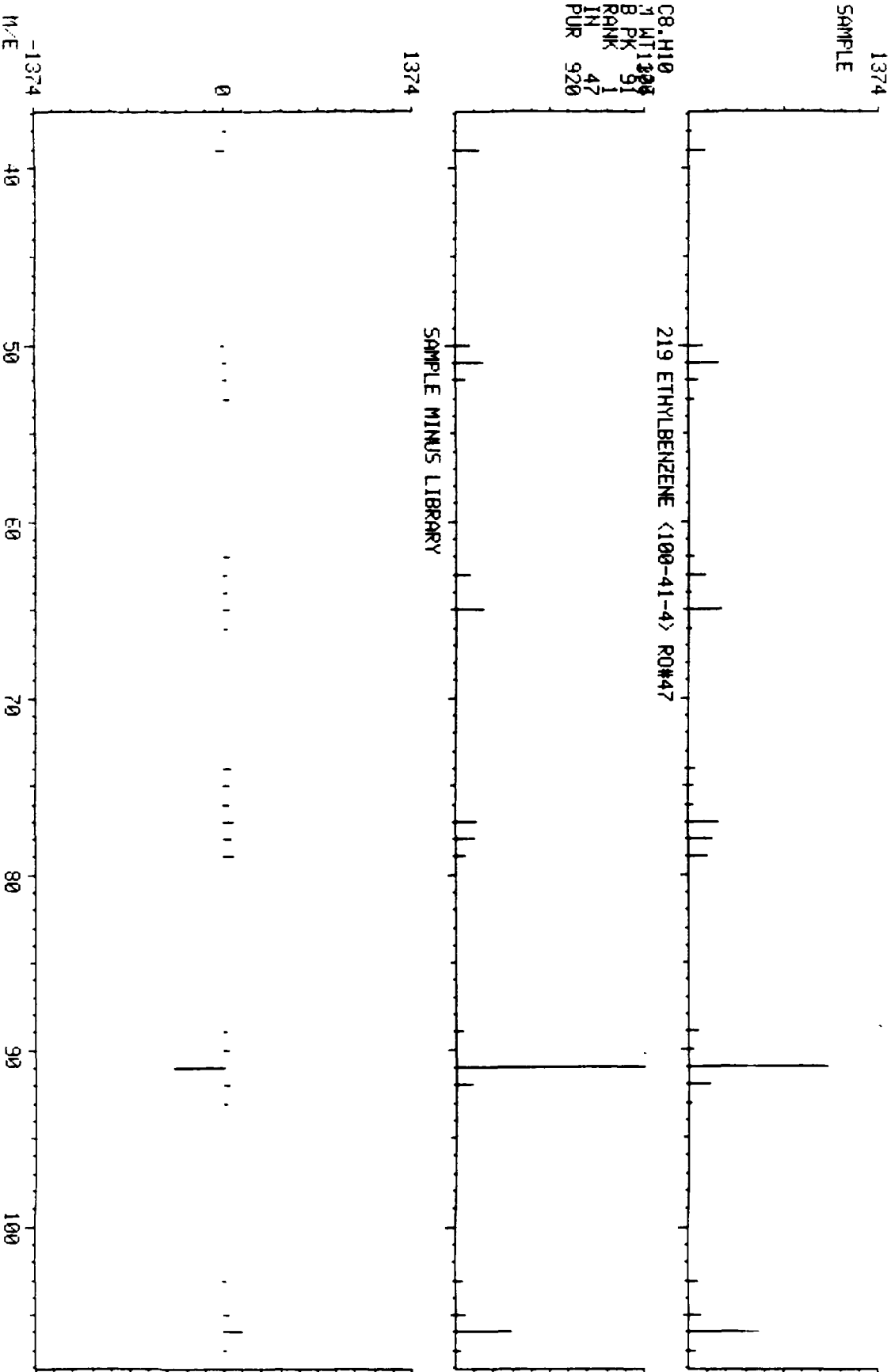
DATA: GM009679C19 #919 BASE M/E: 43/ 43
RIC: 105983.7 140799.



LIBRARY SEARCH
12/22/89 4:18:00 + 12:52
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5.158 2N 0T)

COMFUCHEM LABS
DATA: GH009679C19 #1029
BASE M/E: 91
RIC: 338431.

C8.H10
1 M/I 100
B PK 91
RANK 47
IN 1
PUR 920

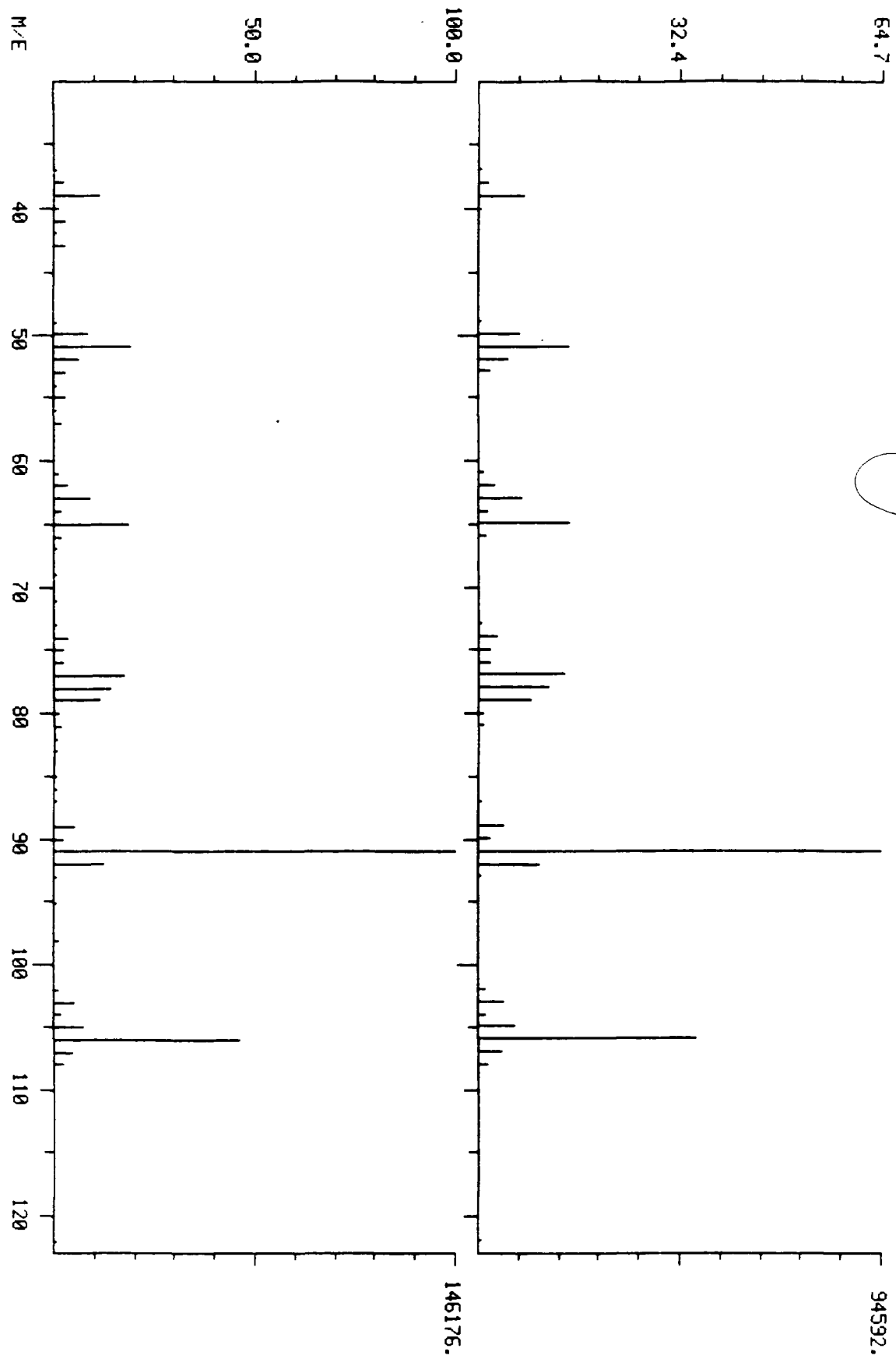


DUAL MASS SPECTRUM
12/22/89 4:18:00 + 12:52
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N)

219 ETHYLBENZENE <100-41-4> R0#47

COMPUCHEM LABS

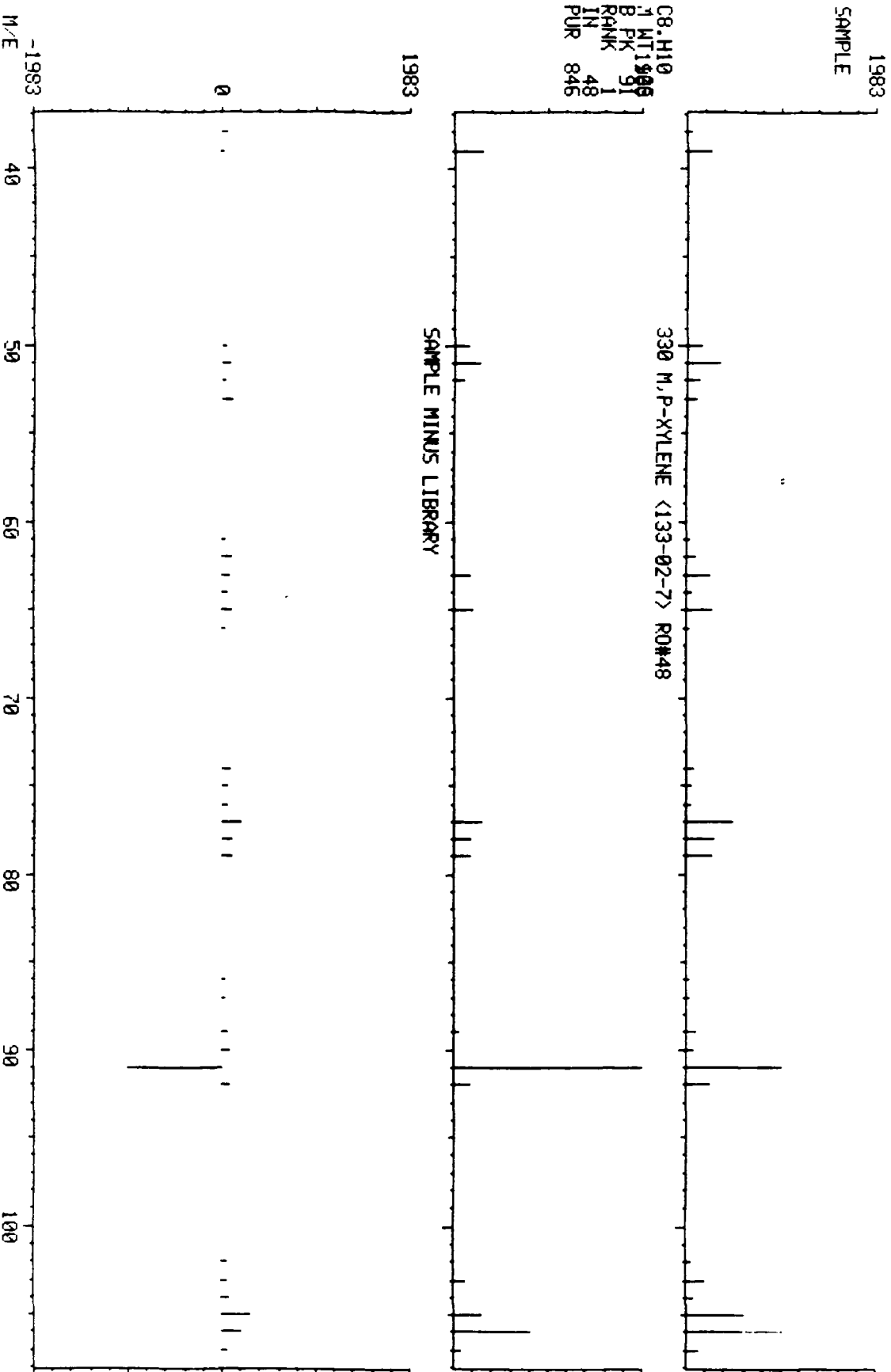
DATA: GH09679C19 #1029 BASE M/E: 91/ 91
RIC: 350207./ 514047.



COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 4:18:00 + 13:07
SAMPLE: 5G CC#809679 CASE#18756.7 EPA#B201A QN#19
ENHANCED (5 158 2N 0T)

DATA: GH009679C19 #1049

BASE M/E: 91
RIC: 584703.

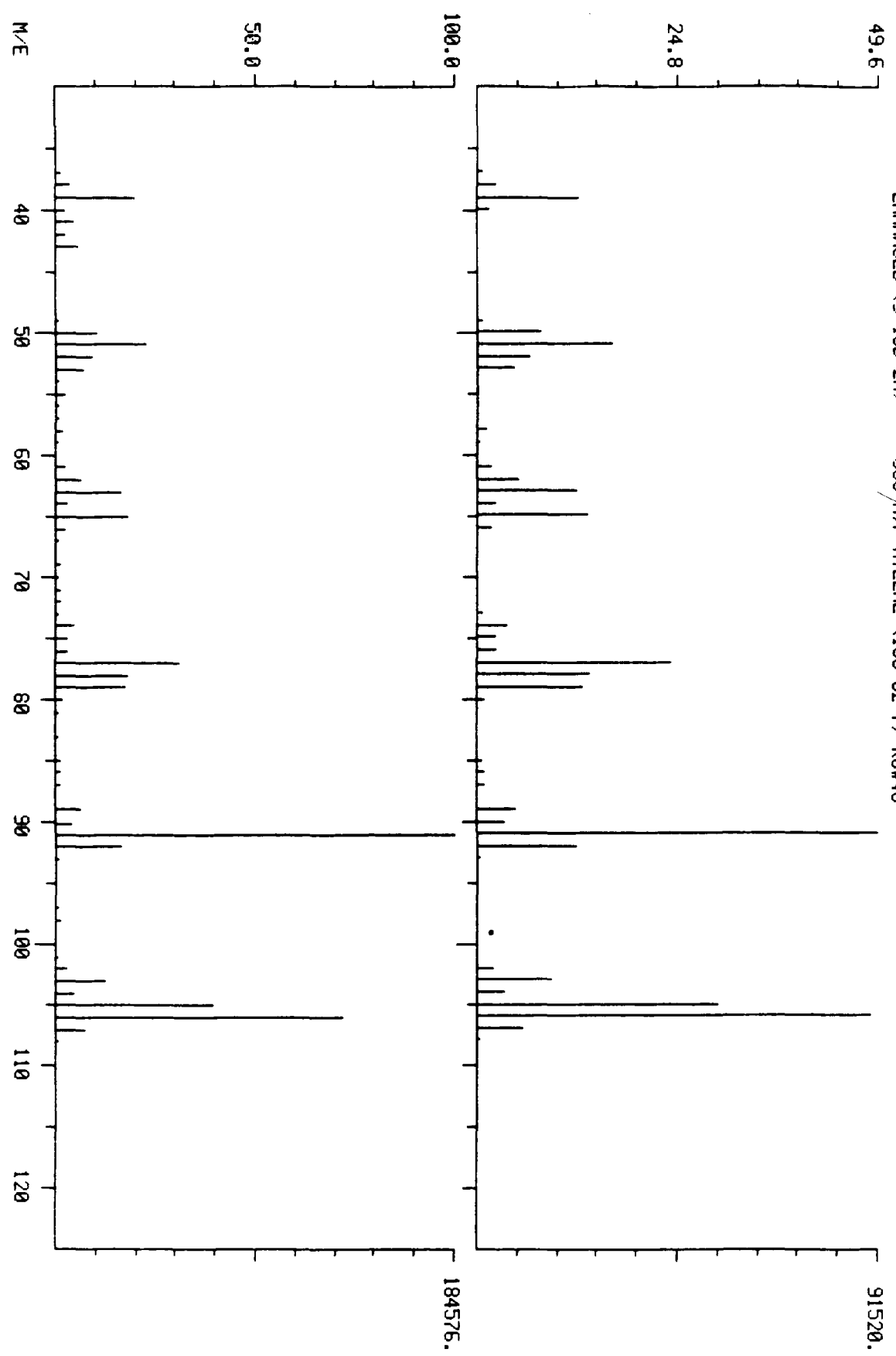


DUAL MASS SPECTRUM
12/22/89 4:18:00 + 13:07
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N) 330 M.P.-XYLENE <133-02-7> R0#48

COMPUCHEM LABS

DATA: GH009679C19 #1049 BASE M/E: 91/ 91

RIC: 595967./ 919551.

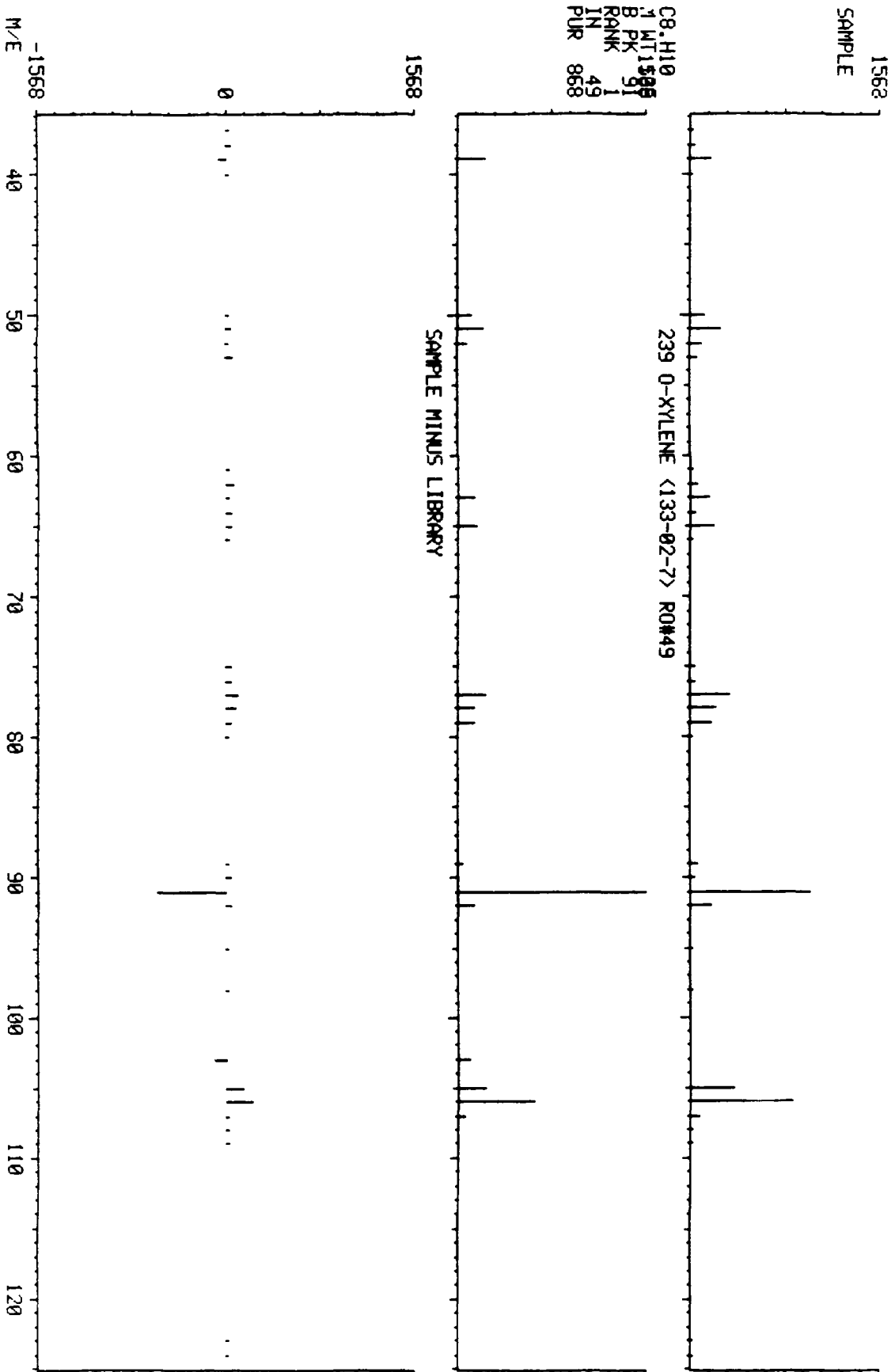


LIBRARY SEARCH
12/22/89 4:18:00 + 13:48
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009679C19 #1104

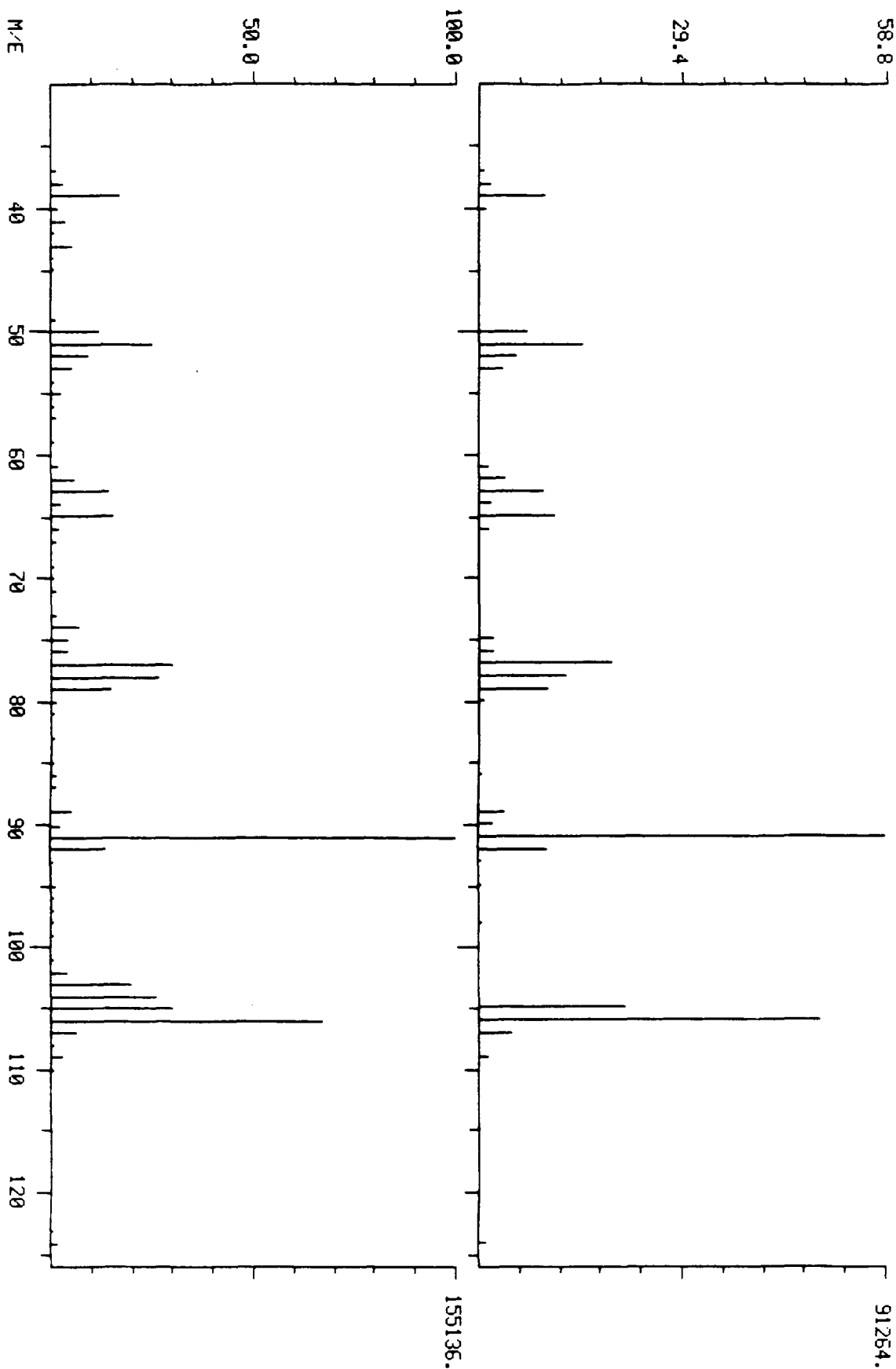
BASE M/E: 91
RIC: 419327.



DUAL MASS SPECTRUM
12/22/89 4:18:00 + 13:48
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 15B 2N) 239 0-XYLENE <133-02-7> R0#49

COMPUCHEM LABS

DATA: GH009679C19 #1104 BASE M/E: 91/ 91
RIC: 420863./ 789503.

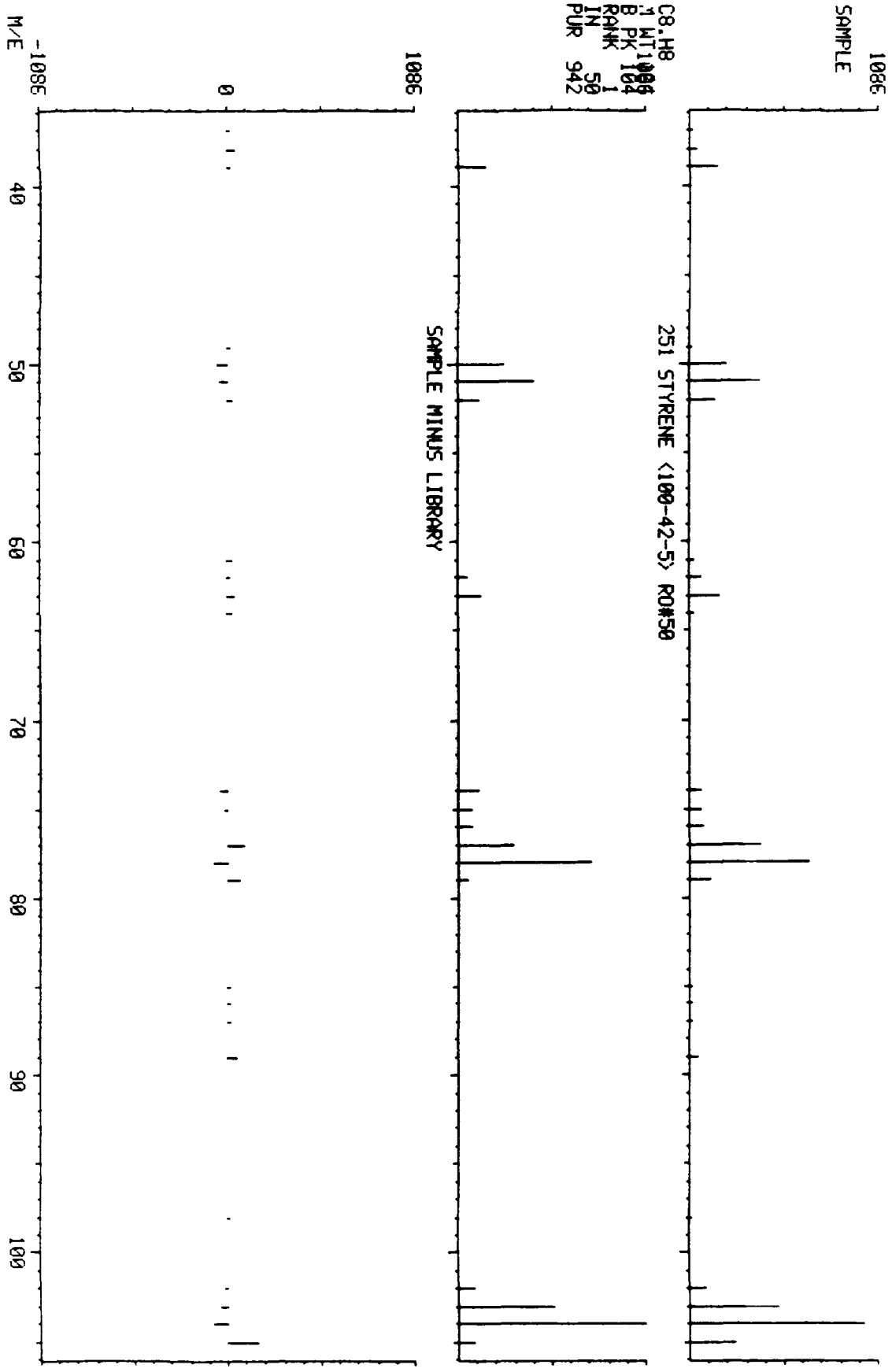


COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 4:18:00 + 13:51
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 15B 2N 0T)

DATA: GH009679C19 #1108

BASE M/E: 104
RIC: 279551.

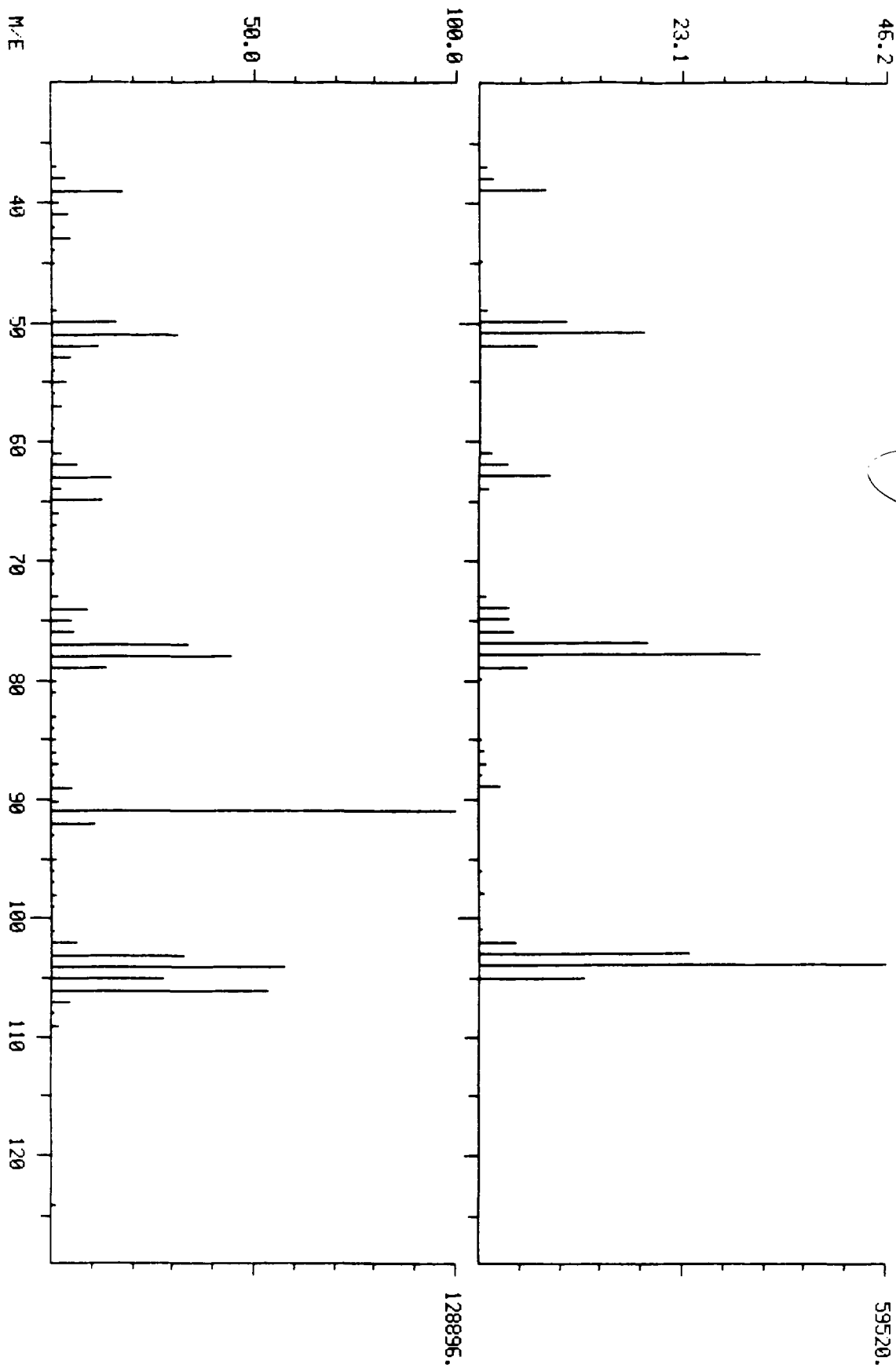
C8.H8
1 WT 1096
B PK 104
RANK 1
IN 50
PUR 942



DUAL MASS SPECTRUM
12/22/89 4:18:00 + 13:51
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 15B 2N) 251 STYRENE <100-42-5> F0#50

COMPUCHEM LABS

DATA: GH009679C19 #1108 BASE M/E: 104/ 91
RIC: 282111.7 740351.



LIBRARY SEARCH
 12/22/89 4:18:00 + 8:13
 SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A QN#19
 ENHANCED (5 158 24 0T)

COMPUchem LABS

DATA: GH009679C19 # 658

BASE M/E: 96
 RIC: 165119.

1042
 SAMPLE

06.H8.0

M WT 1042
 B PK 96
 RANK 1
 IN 776
 PUR 943

FURAN, 2,5-DIMETHYL - CAS# 625-86-5

06.H10.02

M WT 1042
 B PK 114
 RANK 67
 IN 1996
 PUR 616

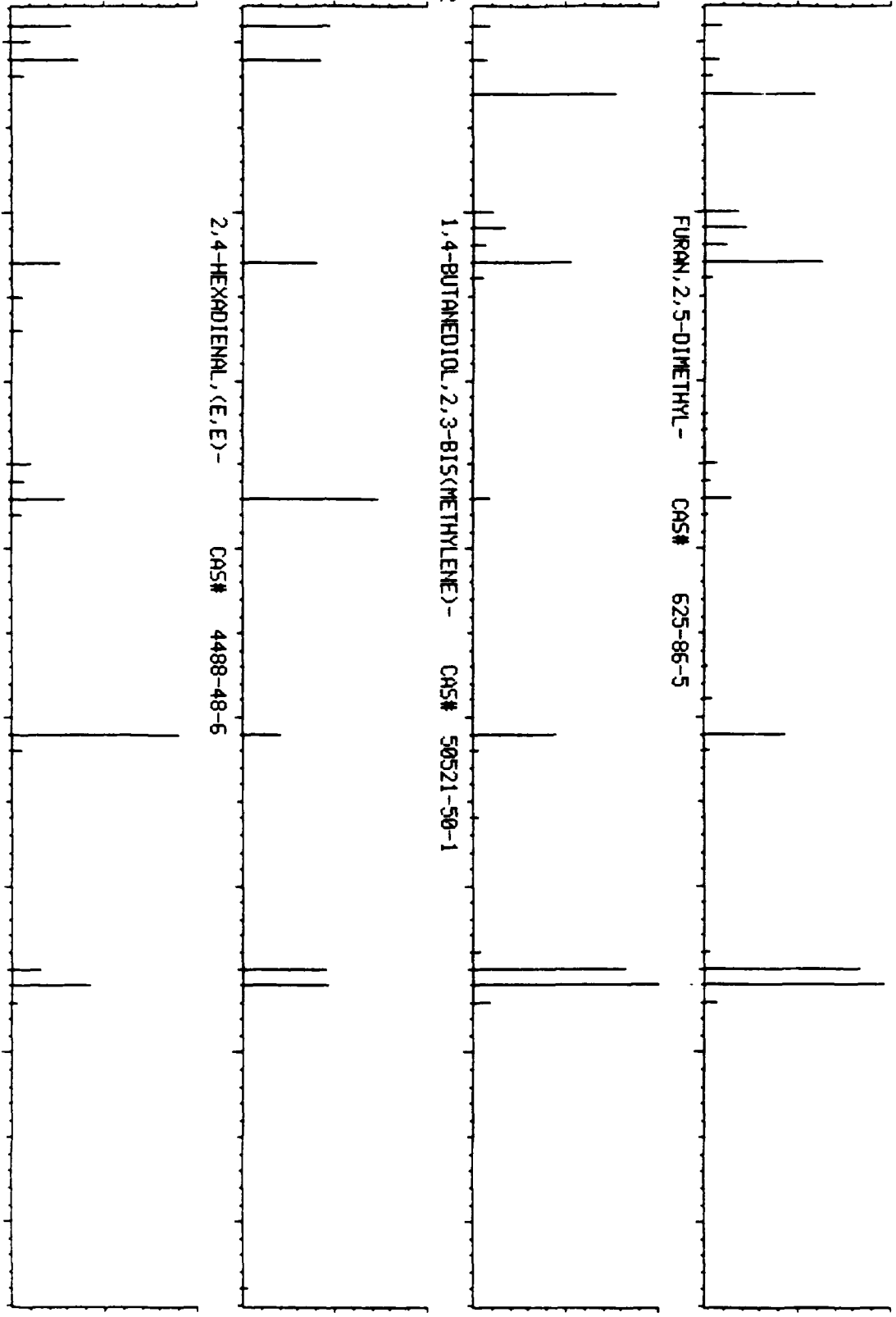
1,4-BUTANEDIOL, 2,3-BIS(METHYLENE) - CAS# 50521-50-1

06.H8.0

M WT 1042
 B PK 96
 RANK 81
 IN 781
 PUR 607

2,4-HEXADIENAL, (E,E) - CAS# 4488-48-6

M/E 40 50 60 70 80 90 100 110



LIBRARY SEARCH
 12/22/89 4:18:00 + 11:01
 SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
 ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009679C19 # 882

BASE M/E: 95
 RIC: 172799.

1000
 SAMPLE

C6.H6.02
 1000
 M WT 110
 B PK 95
 RANK 1
 IN 1598
 PUR 810

ETHANONE,1-(2-FURANYL)- CAS# 1192-62-7

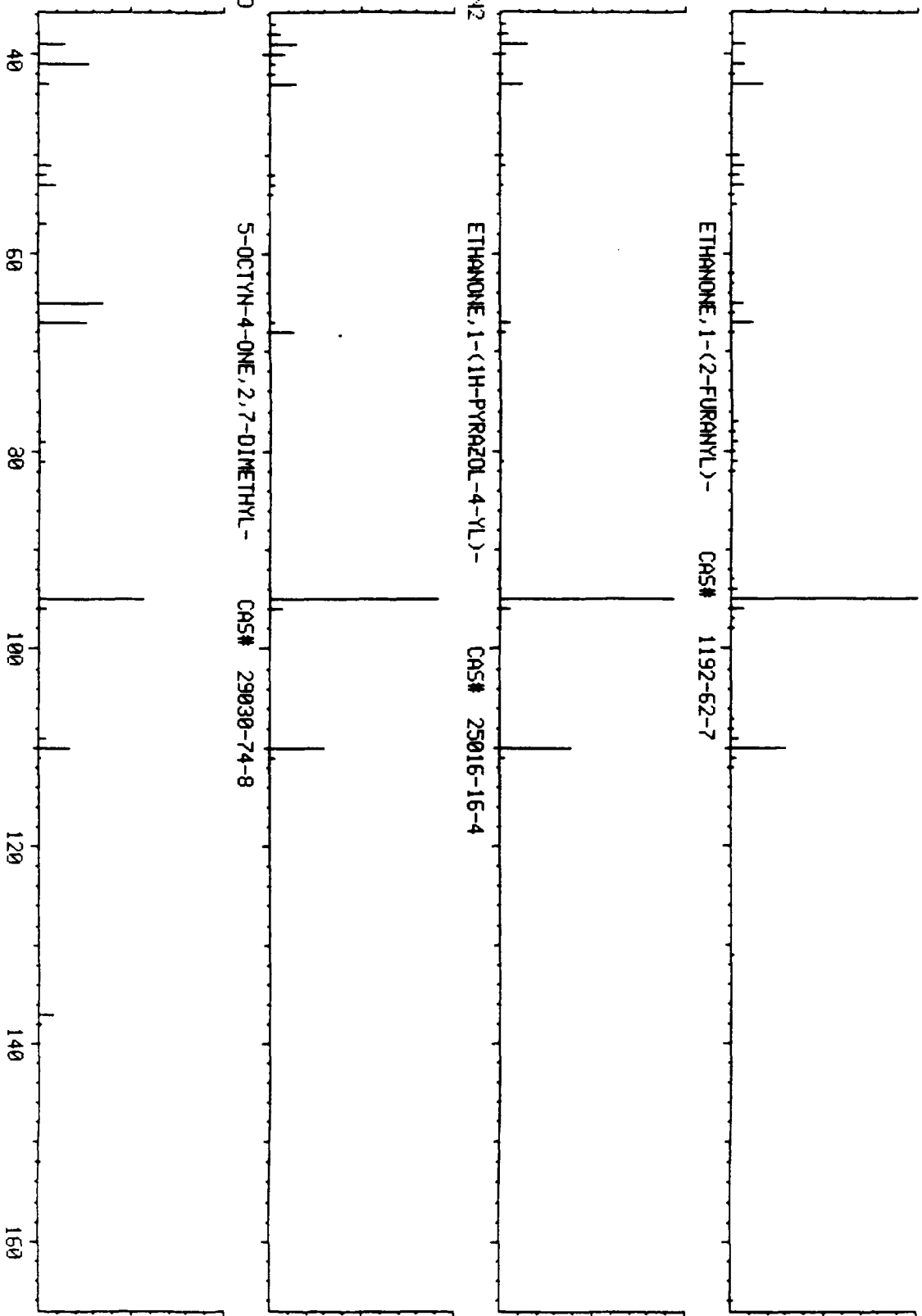
C5.H6.0.N2
 1000
 M WT 110
 B PK 95
 RANK 2
 IN 1588
 PUR 765

ETHANONE,1-(1H-PYRAZOL-4-YL)- CAS# 25016-16-4

C10.H16.0
 1000
 M WT 152
 B PK 95
 RANK 3
 IN 6245
 PUR 728

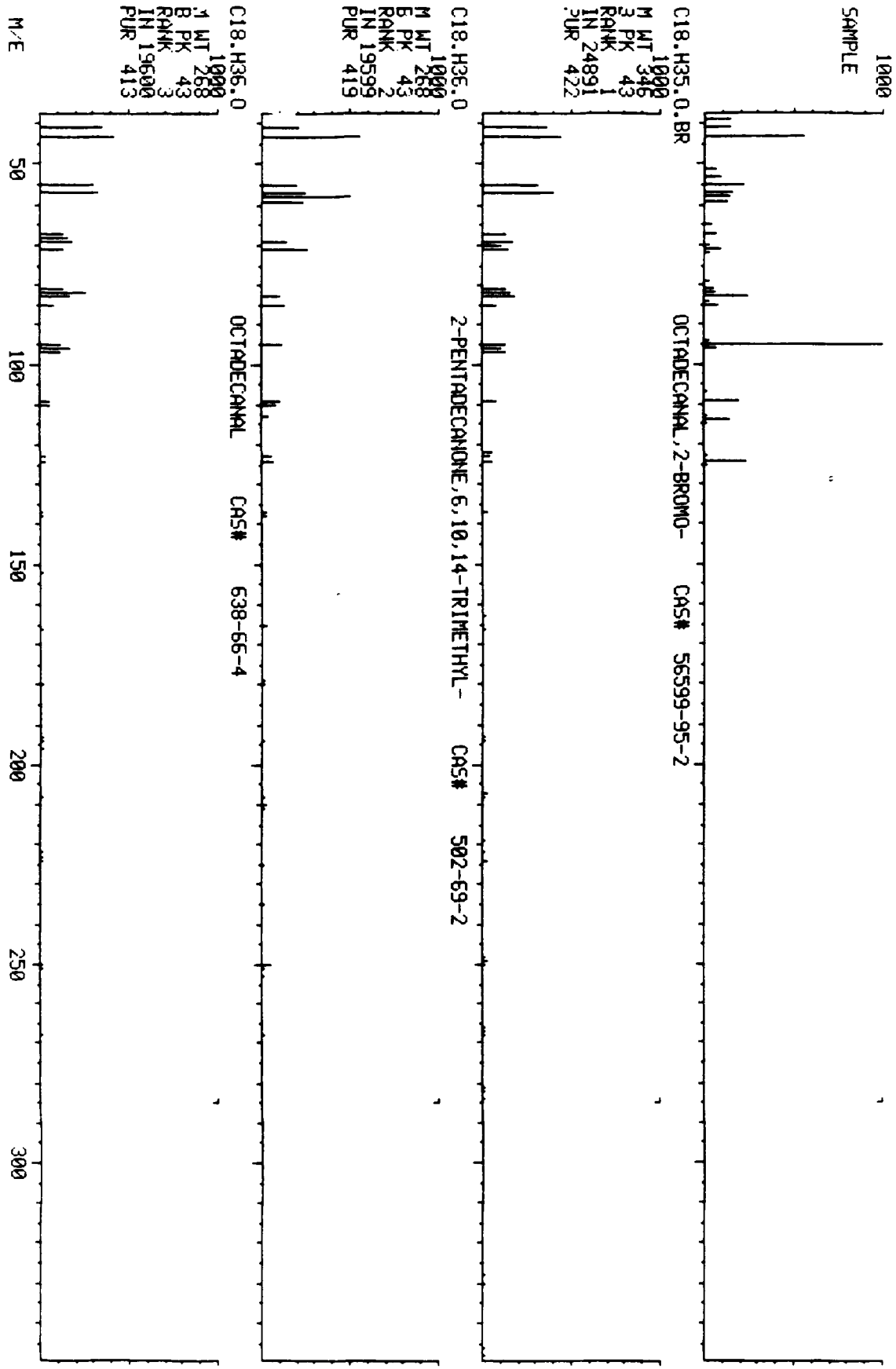
5-OCTYN-4-ONE,2,7-DIMETHYL- CAS# 29030-74-8

M/E 40 50 60 80 100 120 140 160



LIBRARY SEARCH
12/22/89 4:18:00 + 13:25
SAMPLE: 5G CC#809679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (S 158 2N 0T)

COMPUchem LABS
DATA: GH009679C19 #1075
BASE M/E: 95
RIC: 156927.

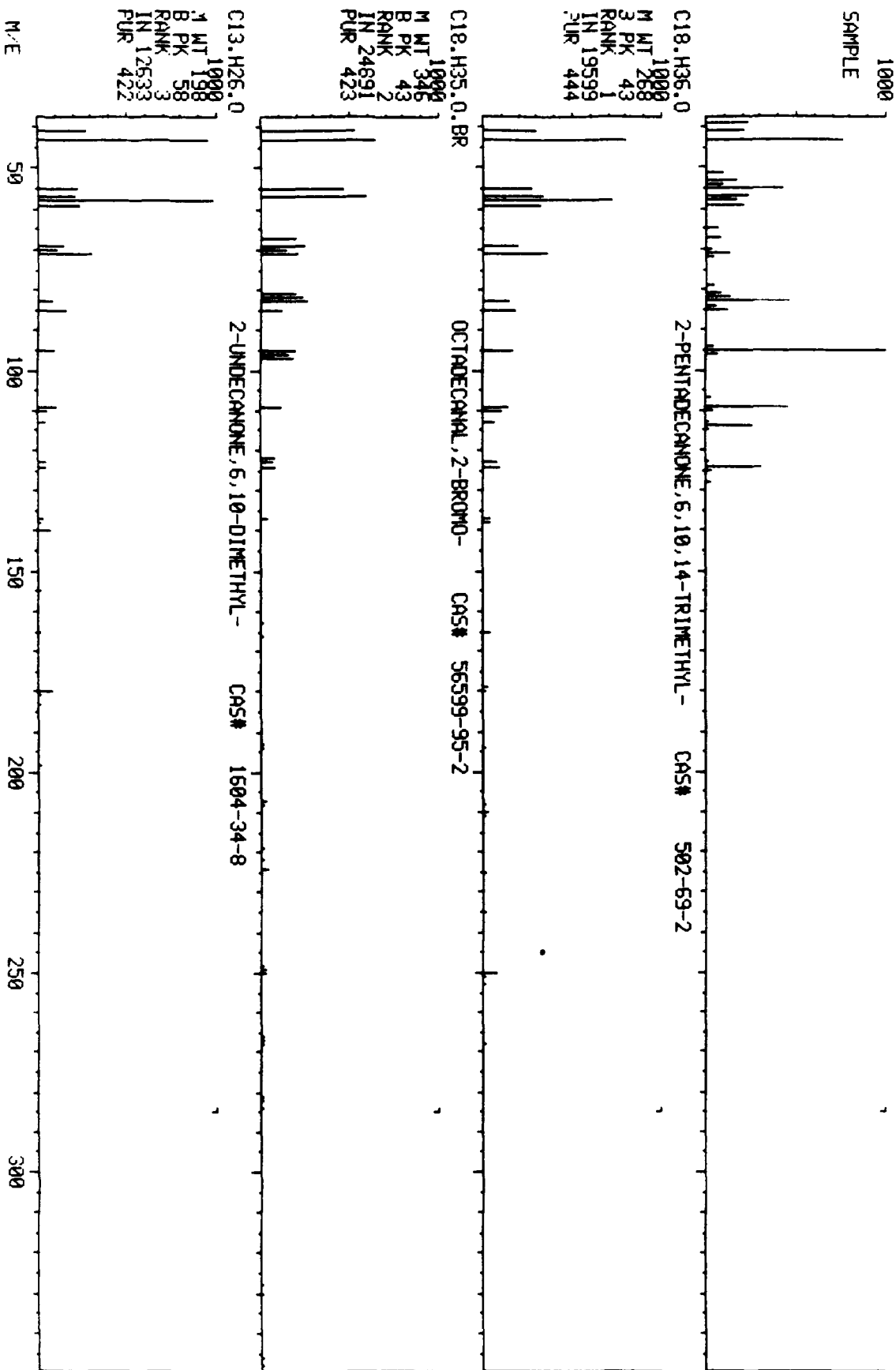


LIBRARY SEARCH
12/22/89 4:18:00 + 13:27
SAMPLE: SG CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009679C19 #1076

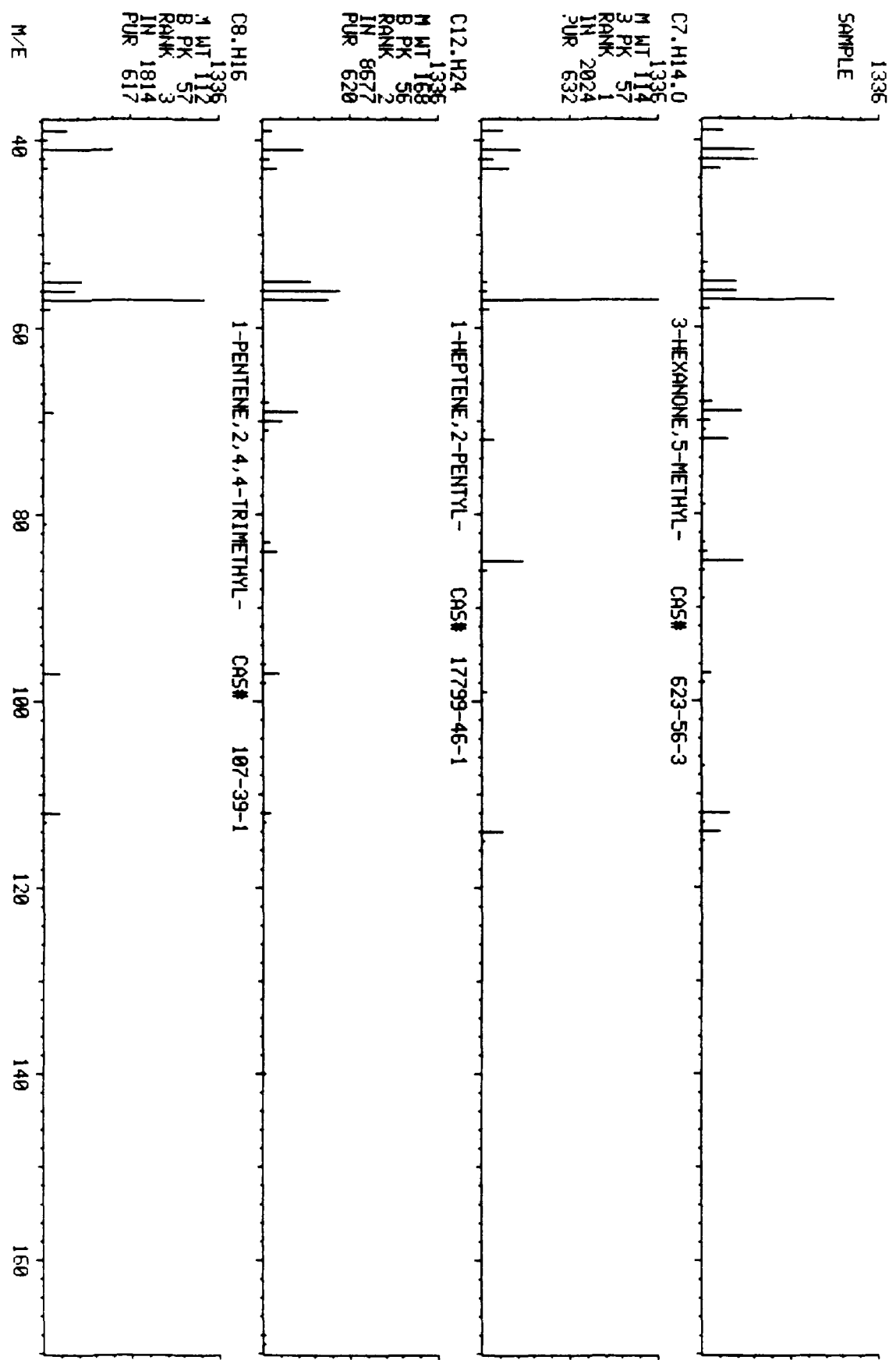
BASE M/E: 95
RIC: 154623.



LIBRARY SEARCH
12/22/89 4:18:00 + 14:04
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (S 158 ZN 01)

COMPUCHEM LABS

DATA: GH009679C19 #1125
BASE M/E: 57
RIC: 169215.



LIBRARY SEARCH
 12/22/89 4:18:00 + 14:33
 SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
 ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

DATA: CH009579C19 #1164

BASE M/E: 105
 RIC: 72191.

1028
 SAMPLE

C9.H12
 1028
 M WT 120
 3 PK 105
 RANK 1
 IN 2524
 PUR 592

BENZENE, (1-METHYLETHYL)- CAS# 98-82-8

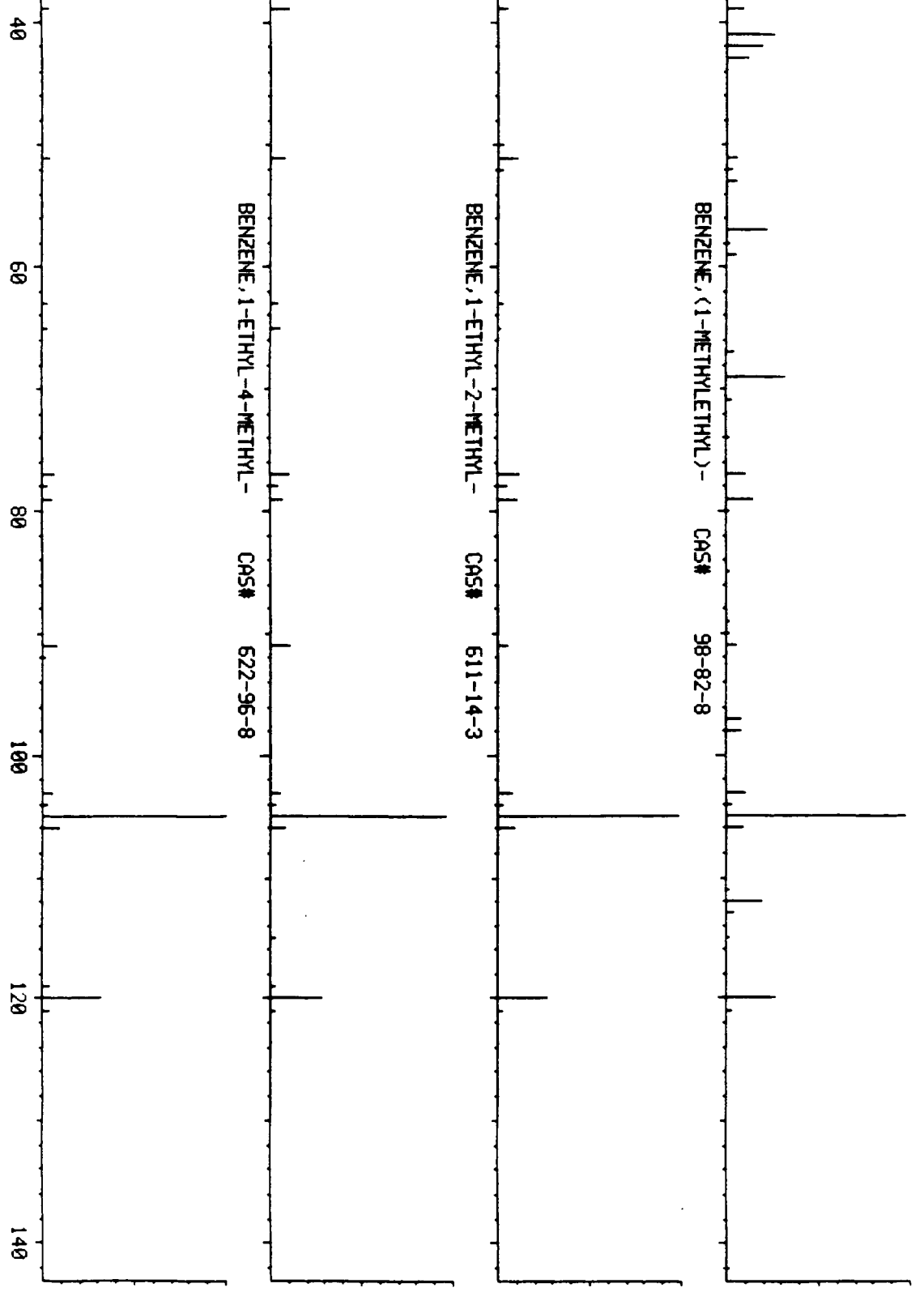
C9.H12
 1028
 M WT 120
 B PK 105
 RANK 2
 IN 2528
 PUR 575

BENZENE, 1-ETHYL-2-METHYL- CAS# 611-14-3

C9.H12
 1028
 1 WT 120
 B PK 105
 RANK 3
 IN 2530
 PUR 570

BENZENE, 1-ETHYL-4-METHYL- CAS# 622-96-8

M/E

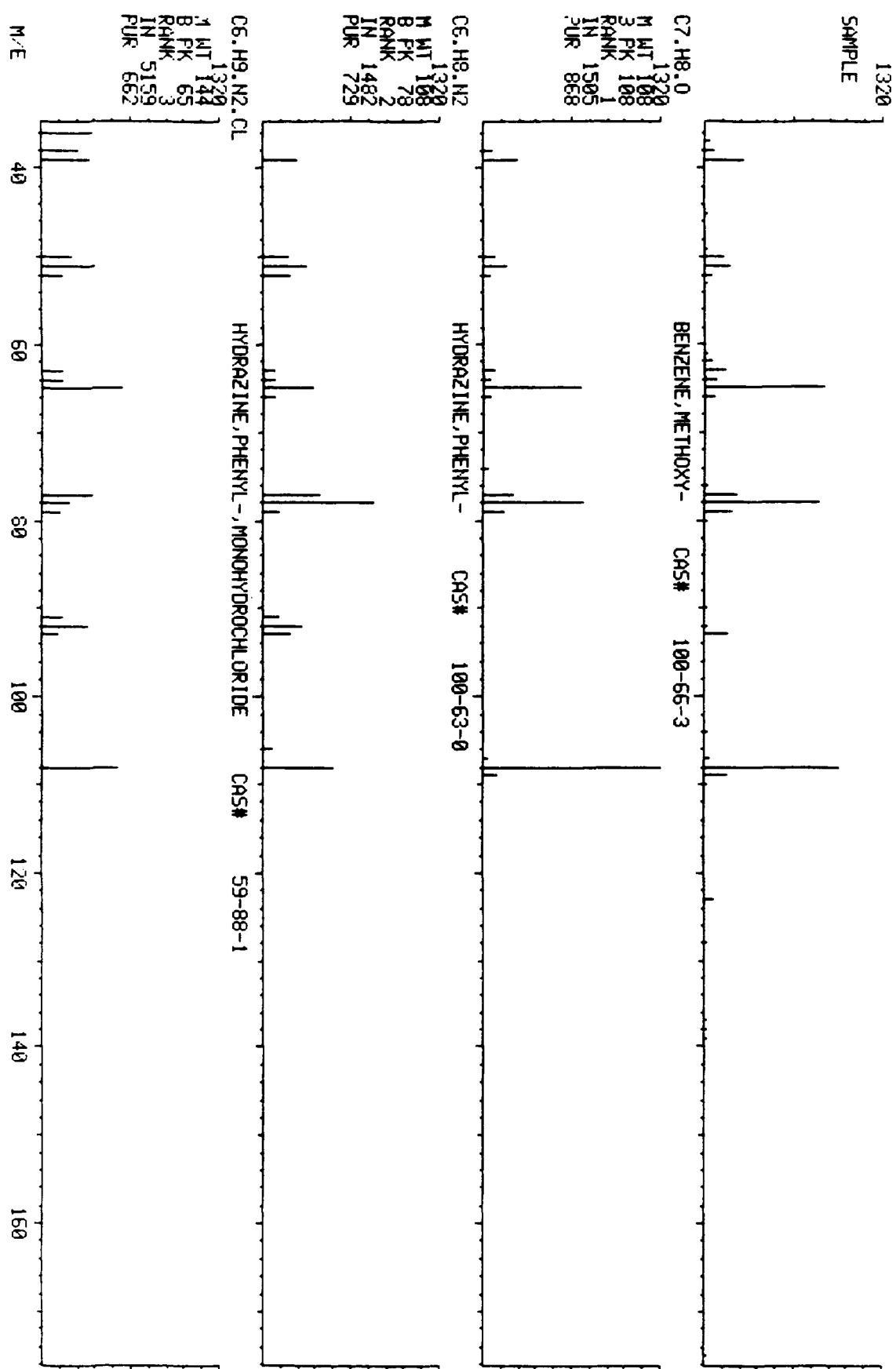


LIBRARY SEARCH
 12/22/89 4:18:00 + 14:43
 SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
 ENHANCED (5 15B 2N 0T)

COMPUCHEN LABS

DATA: GH009679C13 #1177

BASE M/E: 108
 RIC: 270847.

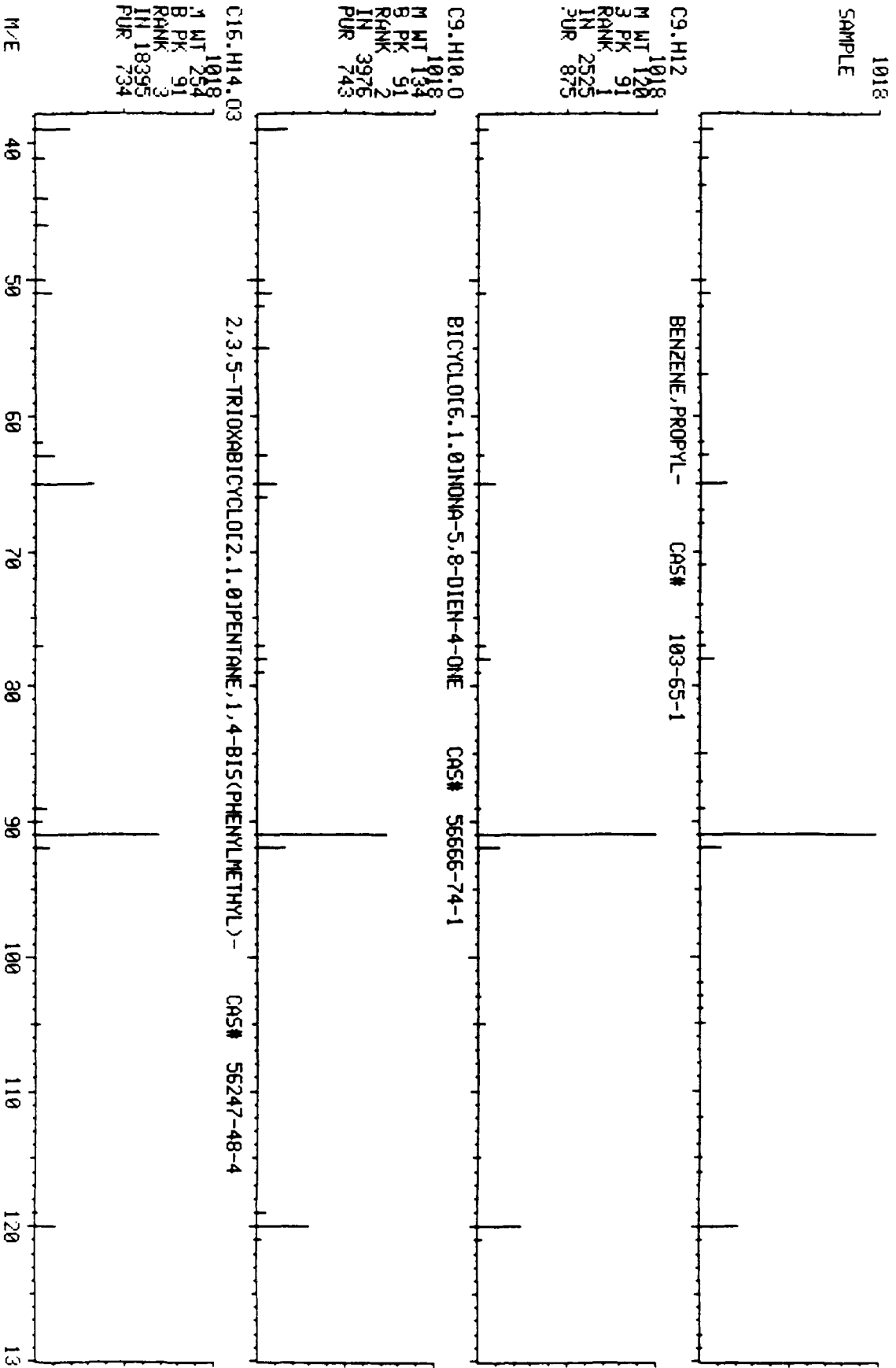


LIBRARY SEARCH
12/22/89 4:18:00 + 15:20
SAMPLE: 5G CC#309679 CASE#19756.7 EPA#B201A GN#19
ENHANCED (5 15B 2N 0T)

COMPUCHEM LABS

DATA: GH009679C19 #1227

BASE M/E: 91
RID: 174079.



LIBRARY SEARCH
12/22/89 4:18:00 + 15:36
SAMPLE: 5G.CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009679C19 #1248

BASE M/E: 105
RIC: 347135.

2015
SAMPLE

C9.H12

M WT 2015
B PK 105
RANK 1
IN 2539
PUR 618

2,3-HEPTADIEN-5-YNE, 2,4-DIMETHYL-

CAS# 41898-89-9

C9.H12

M WT 2015
B PK 105
RANK 2
IN 2524
PUR 572

BENZENE, (1-METHYLETHYL)-

CAS# 98-82-8

C9.H12

M WT 2015
B PK 120
RANK 3
IN 2531
PUR 572

CYCLOHEXENE, 1-(1-PROPYNYL)-

CAS# 1655-05-6

M/E

40 50 60 70 80 90 100 110 120

LIBRARY SEARCH
12/22/89 4:18:00 + 15:04
SAMPLE: 5G CC#309679 CASE#18756.7 EPA#B201A ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009679C19 #1285

BASE M/E: 105
RIC: 323583.

1339
SAMPLE

C9.H12

M WT 1339
B PK 105
RANK 1
IN 2528
PUR 810

BENZENE, 1-ETHYL-2-METHYL-

CAS#

611-14-3

C9.H12

M WT 1339
B PK 105
RANK 2
IN 2529
PUR 803

BENZENE, 1-ETHYL-3-METHYL-

CAS#

620-14-4

C9.H12

M WT 1339
B PK 105
RANK 3
IN 2539
PUR 802

2,3-HEPTADIEN-5-YNE, 2,4-DIMETHYL-

CAS# 41898-89-9

M/E

40

60

80

100

120

140

LAB INSTRUCTIONS:

USE FOR QC!!! SAMPLED DATE 12/18/89

RECEIPT DATE 12/20/89

CASE#: 18756 7

DUE DATE

VOA
GC/MS WORKSHEET

COMPUCHEM#: 309679

R1 [] R2 [] D1 [] ()
R3 [] R4 [] D2 [] ()

L.L. SOLID, EPA SOW 2/88

Sample Prep Code---155
Instrument Code---413
Compound List-----494
Surrogate Std-----394
Internal Std-----036

=====

SAMPLE ID#: B201A Dry Wt. Factor 1.10 % Moisture 9

=====

GC/MS ANALYSIS

Amount Purged: [] 10mls/Xg soil or [] Dilution _____ ul/10000ul/Xg soil
Internal Standard Volume Added 5 ul
Surrogate Standard Volume Added 5 ul
BFB Filename 6589127219 Disk (28574)
Blank Filename 64004795619 Disk ()
Standard Filename 6589127219 Disk ()
Sample Filename 64009679619 Disk ()

ANALYST(S): Injection 1422-16 Work-up 1422-16

GC/MS REVIEW

CONDITION CODE

[]
[DF]

Entry Codes OK, EA, ES, SM, JS, SL, SH, JA, DA

Non-Entry Codes IM, IL, IH, SW, CT, CS, PC, NR
IF, LA, DI, CO, RN, DW, SI, SF
UP, BB, OT, VC, FO, NS

Disposition: [] Complete

Extraneous Peak Search Results:

of Peaks Found: 10

[] Reprep neat required

[] Reprep using 1.0 g

Quality Assurance Notice(s):

Notices Required _____

[] Dilute ()



COMMENTS:

GC/MS Review Sawell Date 1/4/90 Auditor _____ Date _____

REPORT INTEGRATION

Final Reportable Package(s): 64009679619 Total # of Injections: 123209679619

QA COMMENTS:

Initials _____ Date _____

FINAL REVIEW

Initials _____ Date _____

AC1007 (05/89)

VOLATILE PREPARATION WORKSHEET

QUEUE # 19 PREP CODE -155 ASSIGNED TO Harshad Jishi DATE 12/20/89

SAMPLE NUMBER	CASE NUMBER	SDG	QC SAMPLE		SAMPLE WEIGHT (G) VOLUME (ML)	SAMPLE ID	COMMENTS
			TYPE	ORIGINAL			
309763	18520	0164			5.0g	TB-12	
309768	↓	↓			5.0g	WB5-1	
309769	↓	↓			5.0g	WB5-2	
309770	↓	↓			5.0g	WB5-4	
309679	18756	0007			5.0g	B201A	
309680	↓	↓	SS	309679	5.0g	B201A MS	
309681	↓	↓	SS	309679	5.0g	B201A MSD	
309682	↓	↓	BS		0.0g		
309686	↓	↓			5.0g	B201B	
309687	↓	↓			5.0g	B202AR	
309688	↓	↓			5.0g	B202A	
309689	↓	↓			5.0g	B202B	
309690	↓	↓			5.0g	B202C	
309784			B1		5.0 ml	B3	
309785			B2		0.0 ml	B4	
309786			B3		0.0 ml	B5	
309787			B4		0.0 ml	B6	
309788			B5		0.0 ml	B7	

SURROGATE # 4 LOT # 1 MANUAL OPERATOR 739 / 516
 AMOUNT _____
 RELINQUISHED BY [Signature] DATE 12/20/89 RECEIVED BY [Signature] DATE 12/20

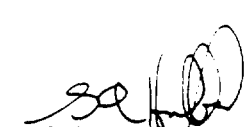
CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	464	34800	50.0		10
221	50	CHLOROMETHANE				BDL	10
231	62	VINYL CHLORIDE				BDL	10
220	94	BROMOMETHANE				BDL	10
209	64	CHLOROETHANE				BDL	10
216	96	1,1-DICHLOROETHENE				BDL	
254	76	CARBON DISULFIDE				BDL	
252	43	ACETONE (2-PROPANONE)			133.0	150 130 B	1
248	114 I	1,4-DIFLUOROBENZENE (IS)	609	134000	50.0		
222	84	METHYLENE CHLORIDE			29.4	32 29 B	
226	96	TRANS-1,2-DICHLOROETHENE				BDL	
214	63	1,1-DICHLOROETHANE				BDL	
257	43	VINYL ACETATE				BDL	10
237	96	CIS-1,2-DICHLOROETHENE				BDL	
253	72	2-BUTANONE			107.0	120 110	10
211	83	CHLOROFORM				BDL	
227	97	1,1,1-TRICHLOROETHANE				BDL	
206	117	CARBON TETRACHLORIDE				BDL	
203	78	BENZENE			233.0	260 230 E	
215	62	1,2-DICHLOROETHANE				BDL	
270	117 I	D5-CHLOROBENZENE (IS) RO#29	1000	158000	50.0		
229	130	TRICHLOROETHENE				BDL	
217	63	1,2-DICHLOROPROPANE				BDL	
212	83	BROMODICHLOROMETHANE				BDL	
218	75	CIS-1,3-DICHLOROPROPENE				BDL	
256	43	4-METHYL-2-PENTANONE			15.5	17 15	10
225	92	TOLUENE			652.0	720 650 E	
250	75	TRANS-1,3-DICHLOROPROPENE				BDL	
228	97	1,1,2-TRICHLOROETHANE				BDL	
224	164	TETRACHLOROETHENE				BDL	
255	43	2-HEXANONE			524.0	550 520 E	10
208	129	DIBROMOCHLOROMETHANE				BDL	
207	112	CHLOROBENZENE				BDL	
219	106	ETHYLBENZENE			383.0	380 420 E	
330	106	M, P-XYLENE			653.0	650 720 E	
239	106	O-XYLENE			520.0	570 520 E	
251	104	STYRENE			190.0	210 190	
205	173	BROMOFORM				BDL	
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	
258	65 S	D4-1,2-DICHLOROETHANE RO#57			48.3	97. %	
247	95 S	BROMOFLUOROBENZENE			45.2	90. %	
233	98 S	D8-TOLUENE RO#59			39.9	80. %	
289	106	XYLENES (TOTAL)			1170.0	1200 E	5

CORRECTED/REVIEWED BY [Signature]
(GC/MS DATA REVIEWER)

DATE 11/3/90

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT. LIMIT (UG/KG)
299	96	1,2-DICHLOROETHENE (TOTAL)				BDL	
CHECKSUMS:							
		3979.	2073	326800.	4893.3		4892.

CORRECTED/REVIEWED BY



(GC/MS DATA REVIEWER)

DATE

11/2/90

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P	F
40	258	D4-1,2-DICHLOROETHANE RO#57	48.3	50.0	97.	70-121	X	
41	247	BROMOFLUOROBENZENE	45.2	50.0	90.	74-121	X	
42	233	D8-TOLUENE RO#59	39.9	50.0	80.	81-117		

* ADVISORY SURROGATE ONLY

++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

CORRECTION FACTOR CALCULATION:

$$\frac{5.0 \text{ G}}{\text{WET WEIGHT OF SAMPLE (G)}} \times \frac{\text{GC/MS DILUTION FACTOR}}{\text{GC/MS DILUTION FACTOR}} \times \text{DRY WEIGHT FACTOR} =$$

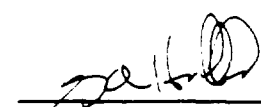
$$\frac{5.0 \text{ G}}{5.00 \text{ (G)}} \times 1.0 \times \frac{1.10}{1.00} = 1.10$$

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:


THE SURROGATES ARE ADDED TO THE SAMPLE PRIOR TO SPARGING.
SURROGATE SPIKE CONVERSION FACTOR = 1.

VERSION 8

CORRECTED/REVIEWED BY


 (GC/MS DATA REVIEWER)

DATE



1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201ARE

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C3R09679B13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 9 Date Analyzed: 12/28/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	(ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	1400	U
74-83-9	-----Bromomethane	1400	U
75-01-4	-----Vinyl Chloride	1400	U
75-00-3	-----Chloroethane	1400	U
75-09-2	-----Methylene Chloride	440	BDJ
67-64-1	-----Acetone	1400	U
75-15-0	-----Carbon Disulfide	690	U
75-35-4	-----1,1-Dichloroethene	690	U
75-34-3	-----1,1-Dichloroethane	690	U
540-59-0	-----1,2-Dichloroethene (total)	690	U
67-66-3	-----Chloroform	690	U
107-06-2	-----1,2-Dichloroethane	690	U
78-93-3	-----2-Butanone	1400	U
71-55-6	-----1,1,1-Trichloroethane	690	U
56-23-5	-----Carbon Tetrachloride	690	U
108-05-4	-----Vinyl Acetate	1400	U
75-27-4	-----Bromodichloromethane	690	U
78-87-5	-----1,2-Dichloropropane	690	U
10061-01-5	-----cis-1,3-Dichloropropene	690	U
79-01-6	-----Trichloroethene	690	U
124-48-1	-----Dibromochloromethane	690	U
79-00-5	-----1,1,2-Trichloroethane	690	U
71-43-2	-----Benzene	630	DJ
10061-02-6	-----Trans-1,3-Dichloropropene	690	U
75-25-2	-----Bromoform	690	U
108-10-1	-----4-Methyl-2-Pentanone	1400	U
591-78-6	-----2-Hexanone	630	DJ
127-18-4	-----Tetrachloroethene	690	U
79-34-5	-----1,1,2,2-Tetrachloroethane	690	U
108-88-3	-----Toluene	4200	D
108-90-7	-----Chlorobenzene	690	U
100-41-4	-----Ethylbenzene	3100	D
100-42-5	-----Styrene	1300	D
1330-20-7	-----Total Xylenes	9500	D

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201ARE

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309679
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C3R09679B13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 9 Date Analyzed: 12/28/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 9 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

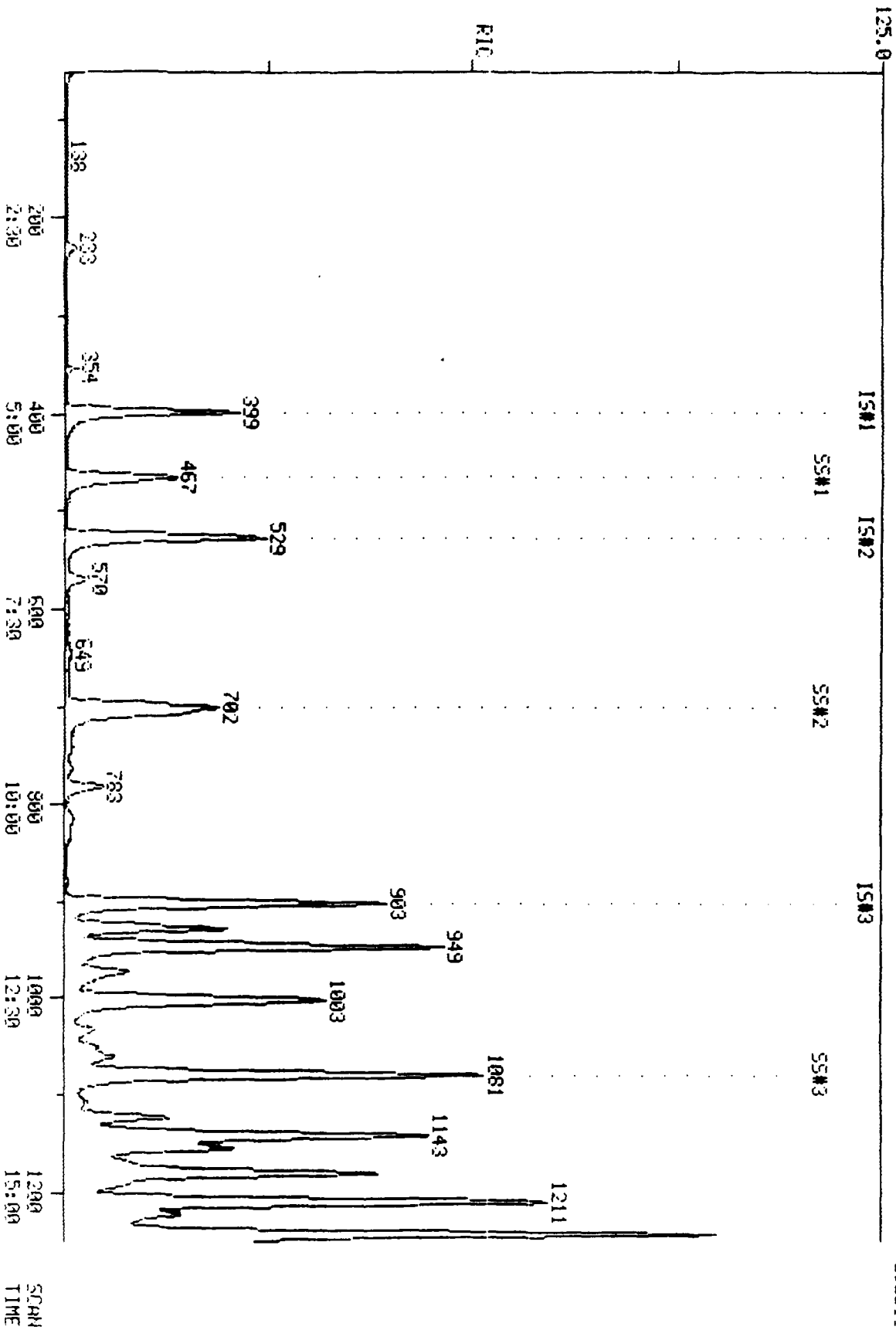
CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 1192-62-7	ETHANONE, 1-(2-FURANYL)-	9.80	820	J
2.	UNKNOWN	12.19	1200	J
3.	SUBSTITUTED BENZENE	14.05	1400	J
4.	SUBSTITUTED BENZENE	14.29	8400	J
5.	SUBSTITUTED BENZENE	14.45	2300	J
6.	SUBSTITUTED BENZENE	14.77	7000	J
7.	SUBSTITUTED BENZENE	15.14	9900	J
8. 496-11-7	1H-INDENE, 2,3-DIHYDRO-	15.30	960	J
9. 271-89-6	BENZOFURAN	15.55	9200	J

RIC
 12/28/83 20:34:00
 SAMPLE: 100UL CC#309579 EPA#B291A CASE#18756 ON #13
 COND5.:

COMPUchem LABS

COMPUchem DATA: C909579R13 SCANS 53 TO 1250

232960.



QUANTITATION REPORT FILE: C3R09679B13
 DATA: C3R09679B13.TI
 12/28/89 20:34:00
 SAMPLE: 100UL CC#309679 EPA#B201A CASE#18756 ON #13
 CONDS.:
 SUBMITTED BY: 13 ANALYST: 1452

RE sat 1/4/90

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
 RESP. FAC. FROM LIBRARY ENTRY

- NO NAME
- 1 *234 BROMOCHLOROMETHANE (IS) <75-97-5> NA#1
- 2 221 CHLOROMETHANE <74-87-3> NA#2
- 3 231 VINYL CHLORIDE <75-01-4> NA#3
- 4 220 BROMOMETHANE <78-83-9> NA#4
- 5 209 CHLOROETHANE <75-00-3> NA#5
- 6 216 1,1-DICHLOROETHENE <75-35-4> NA#8
- 7 254 CARBON DISULFIDE <75-15-0> NA#9
- 8 252 ACETONE (2-PROPANONE) <67-64-1> NA#13
- 9 *248 1,4-DIFLUOROBENZENE (IS) <540-36-3> NA#14
- 10 222 METHYLENE CHLORIDE <75-09-2> NA#16
- 11 226 TRANS-1,2-DICHLOROETHENE <156-60-5> NA#17
- 12 214 1,1-DICHLOROETHANE <75-34-3> NA#19
- 13 257 VINYL ACETATE <108-05-4> NA#20
- 14 237 CIS-1,2-DICHLOROETHENE <156-59-2> NA#21
- 15 253 2-BUTANONE <78-93-3> NA#22
- 16 211 CHLOROFORM <67-66-2> NA#23
- 17 227 1,1,1-TRICHLOROETHANE <71-55-6> NA#24
- 18 206 CARBON TETRACHLORIDE <56-23-5> NA#25
- 19 203 BENZENE <71-43-2> NA#26
- 20 215 1,2-DICHLOROETHANE <107-06-2> NA#27
- 21 *270 D5-CHLOROENZENE (IS) <XX-XX-X> NA#29
- 22 229 TRICHLOROETHENE <79-01-6> NA#30
- 23 217 1,2-DICHLOROPROPANE <78-87-5> NA#31
- 24 212 BROMODICHLOROMETHANE <75-27-4> NA#33
- 25 218 CIS-1,3-DICHLOROPROPENE <10061-1-5> NA#35
- 26 256 4-METHYL-2-PENTANONE <108-01-1> NA#36
- 27 225 TOLUENE <108-88-3> NA#37
- 28 250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> NA#38
- 29 228 1,1,2-TRICHLOROETHANE <79-00-5> NA#39
- 30 224 TETRACHLOROETHENE <127-18-4> NA#41
- 31 255 2-HEXANONE <591-78-6> NA#42
- 32 208 DIBROMOCHLOROMETHANE <124-48-1> NA#43
- 33 207 CHLOROBENZENE <108-90-7> NA#45
- 34 219 ETHYLBENZENE <100-41-4> NA#47
- 35 330 M,P-XYLENE <133-02-7> NA#48
- 36 239 O-XYLENE <133-02-7> NA#49
- 37 251 STYRENE <100-42-5> NA#50
- 38 205 BROMOFORM <75-25-2> NA#51
- 39 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> NA#54
- 40 #258 D4-1,2-DICHLOROETHANE NA#57
- 41 #247 BROMOFLUOROBENZENE <460-00-4> NA#58
- 42 #233 D8-TOLUENE NA#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	399	4:59	1	1.000	A BB	49850.	50.000 UG/L	11.33
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	NOT FOUND							
9	114	529	6:37	9	1.000	A BB	183215.	50.000 UG/L	11.33
10	84	230	2:52	1	0.576	A BB	4693.	3.220 UG/L	0.73
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	NOT FOUND							
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	464	5:48	9	0.877	A BB	14911.	4.601 UG/L	1.04
20	62	NOT FOUND							
21	117	903	11:17	21	1.000	A BB	233794.	50.000 UG/L	11.33
22	130	NOT FOUND							
23	63	NOT FOUND							
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	NOT FOUND							
27	92	709	8:52	21	0.785	A BB	47080.	30.288 UG/L	6.86
28	75	NOT FOUND							
29	97	NOT FOUND							
30	164	NOT FOUND							
31	43	833	10:25	21	0.922	A VB	4890.	4.587 UG/L	1.04
32	129	NOT FOUND							
33	112	NOT FOUND							
34	106	930	11:37	21	1.030	A BV	43049.	22.352 UG/L	5.06
35	106	949	11:52	21	1.051	A VB	133550.	40.297 UG/L	9.13
36	106	1003	12:32	21	1.111	A BB	84236.	29.039 UG/L	6.58
37	104	1009	12:37	21	1.117	A BB	48223.	9.512 UG/L	2.15
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	468	5:51	1	1.173	A BB	96951.	55.464 UG/L	12.57
41	95	1081	13:31	21	1.197	A BB	195806.	46.412 UG/L	10.52
42	98	701	8:46	21	0.776	A BV	114936.	45.629 UG/L	10.34

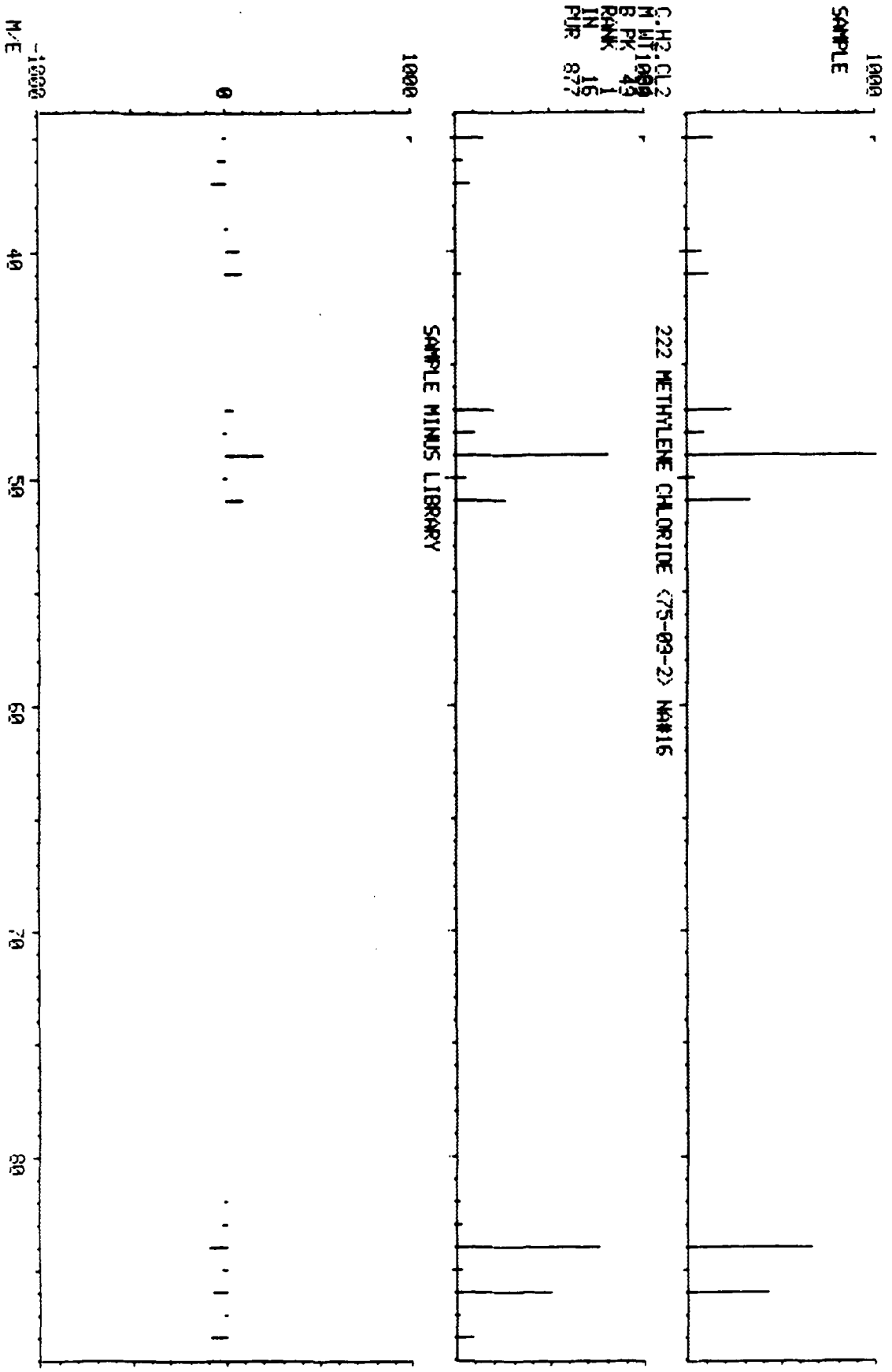
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:07	0.98	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	0:52		10.000			50.00		0.882	
3	0:58		10.000			50.00		1.005	
4	1:11		10.000			50.00		1.237	
5	1:18		10.000			50.00		0.626	
6	2:10		5.000			50.00		1.230	
7	2:16		5.000			50.00		3.765	
8	2:31		10.000			50.00		0.257	
9	6:46	0.98	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	2:58	0.97	5.000	0.12	3.22	50.00	0.094	1.462	0.06
11	3:18		5.000			50.00		1.270	
12	3:56		5.000			50.00		2.282	
13	4:14		10.000			50.00		0.848	
14	4:48		5.000			50.00		1.568	

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:01		10.000			50.00		0.124	
16	5:22		5.000			50.00		2.770	
17	5:24		5.000			50.00		0.713	
18	5:37		5.000			50.00		0.656	
19	5:57	0.97	5.000	0.18	4.60	50.00	0.081	0.884	0.09
20	6:06		5.000			50.00		1.804	
21	11:25	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	7:01		5.000			50.00		0.453	
23	7:22		5.000			50.00		0.350	
24	7:56		5.000			50.00		0.642	
25	8:40		5.000			50.00		0.553	
26	9:07		15.000			50.00		0.233	
27	9:04	0.98	5.000	0.16	30.29	50.00	0.201	0.332	0.61
28	9:40		5.000			50.00		0.243	
29	9:56		5.000			50.00		0.277	
30	9:53		5.000			50.00		0.303	
31	10:35	0.98	15.000	0.06	4.59	50.00	0.021	0.228	0.09
32	10:30		5.000			50.00		0.846	
33	11:28		5.000			50.00		0.966	
34	11:49	0.98	5.000	0.21	22.35	50.00	0.184	0.412	0.45
35	12:04	0.98	5.000	0.21	40.30	50.00	0.571	0.709	0.81
36	12:46	0.98	5.000	0.22	29.04	50.00	0.360	0.620	0.58
37	12:50	0.98	5.000	0.22	9.51	50.00	0.206	1.084	0.19
38	13:06		5.000			50.00		0.762	
39	14:19		5.000			50.00		0.705	
40	6:00	0.98	5.000	0.23	55.46	50.00	1.945	1.753	1.11
41	13:44	0.98	5.000	0.24	46.41	50.00	0.838	0.902	0.93
42	8:57	0.98	5.000	0.16	45.63	50.00	0.492	0.539	0.91

LIBRARY SEARCH
 12/28/89 20:34:00 + 2:52
 SAMPLE: 100UL CC#309679 EPA#B201A CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

COMPUCHEM LABS
 DATA: C3R09679B13 # 230
 BASE M/E: 49
 RIC: 1859.

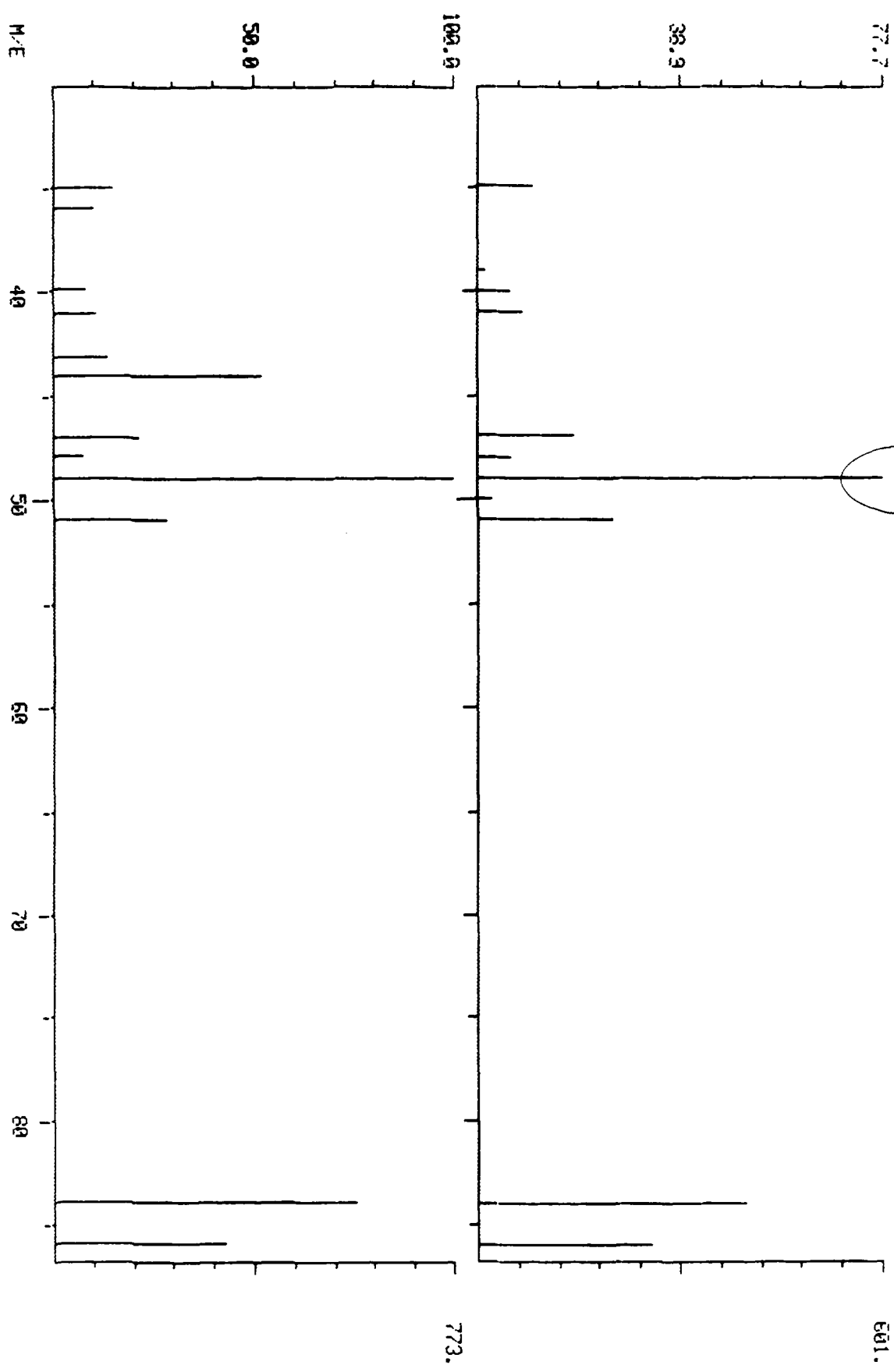
C.H2-CL2
 M WT 1000
 B PK 49
 RANK 1
 IN 15
 PUR 877



DUAL MASS SPECTRUM
12/28/89 20:34:00 + 2:52
SAMPLE: 100UL CC#309679
ENHANCED (S 1SR 2N)

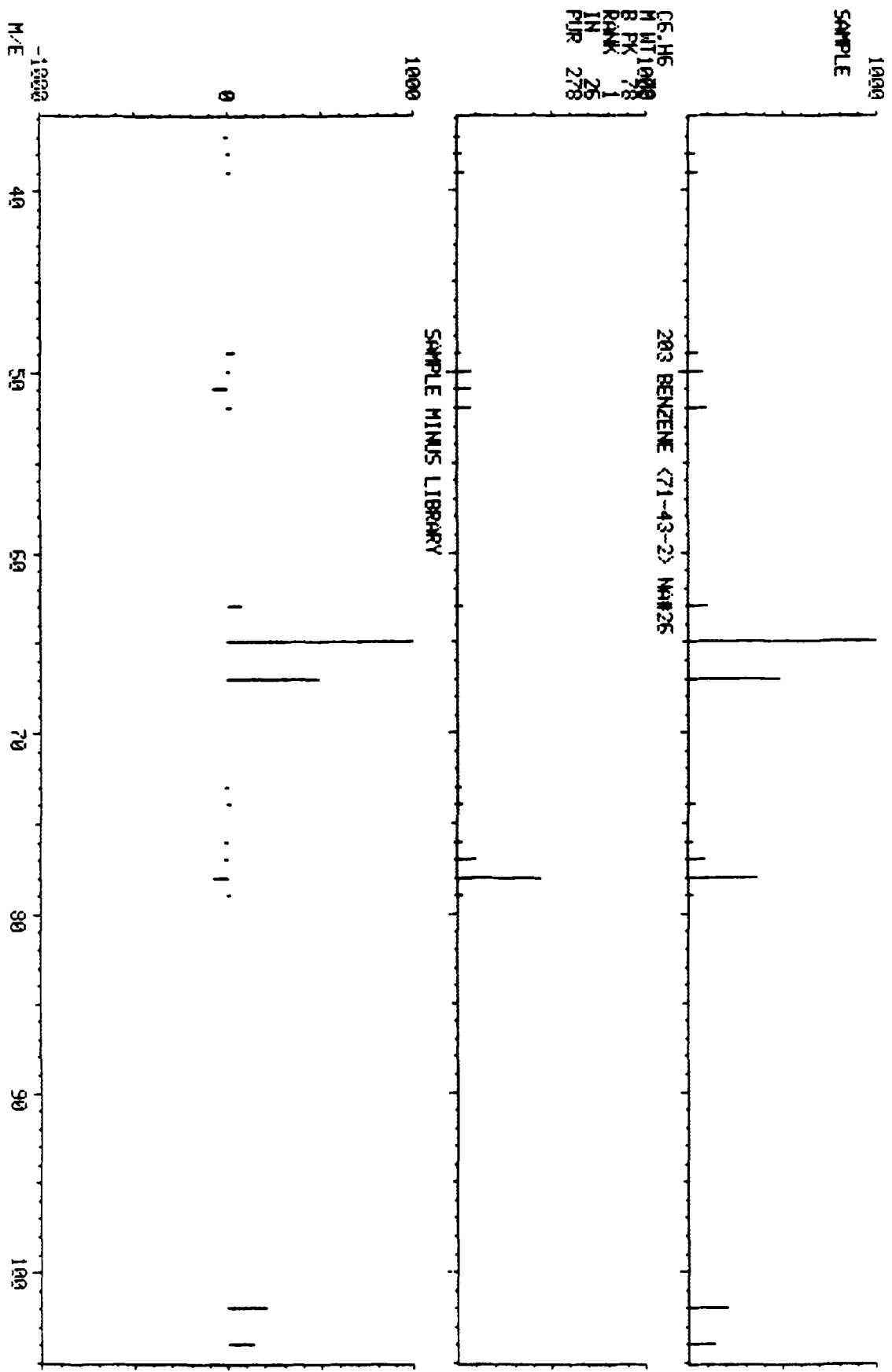
COMPUchem LABS
RE 2004.1.140
EPA#B201A/CASE#18756 ON #13
222 METHYLENE CHLORIDE (75-09-2) NA#15

DATA: C3R09679B13 #230
BASE M/E: 49/
RIC: 1869.7 2959.



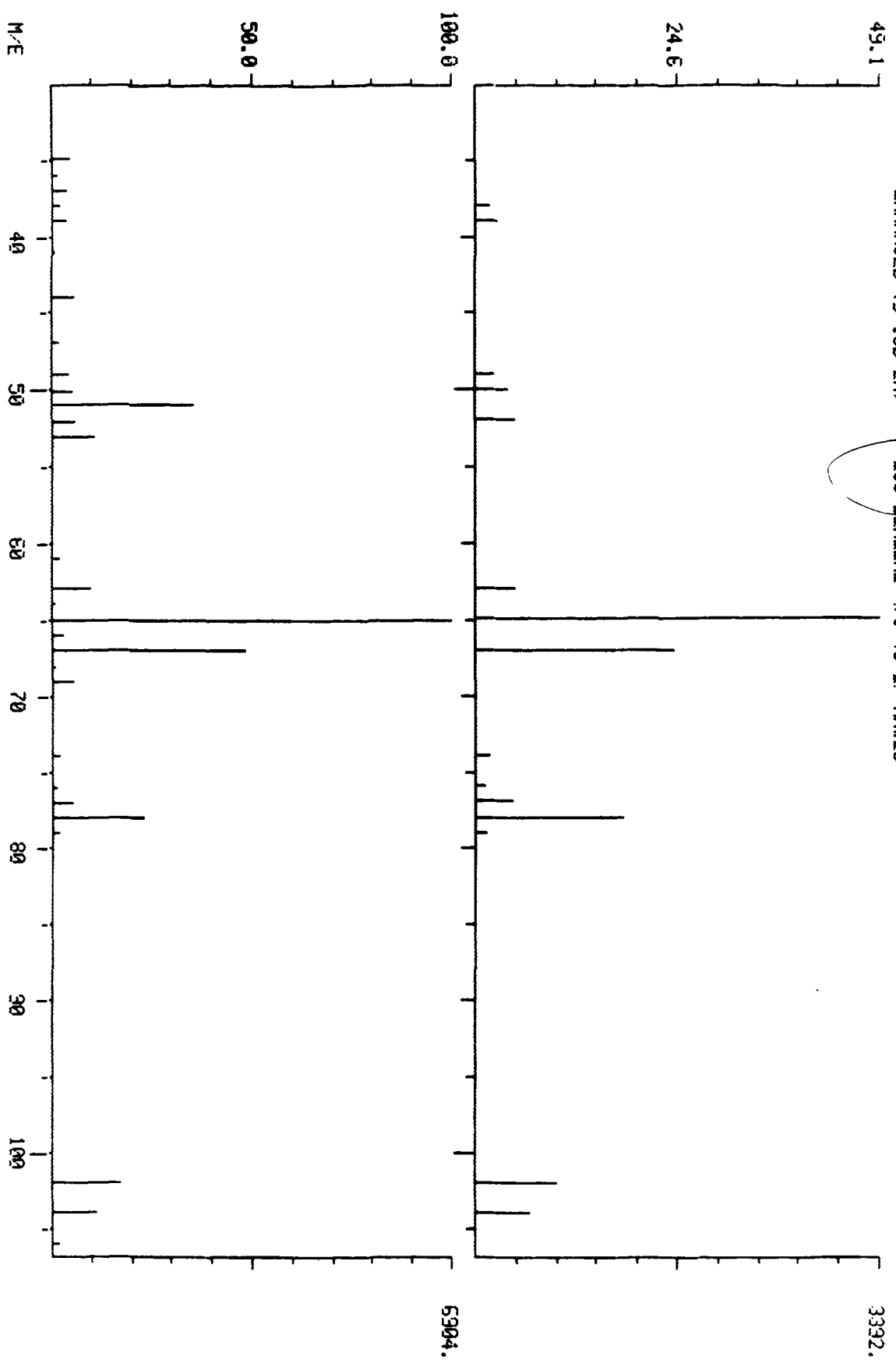
COMPUchem LABS
 LIBRARY SEARCH
 12/28/89 20:34:00 + 5:48 RE *AKM 1/1/90* DATA: C3P09673813 # 464 BASE M/E: 65
 SAMPLE: 100UL CC#309679 EPA#8201A CASE#18756 ON #13 RIC: 9375.
 ENHANCED (5 158 2H 0T)

C6.H6
 M WT 1000
 B PK 78
 RANK 1
 IN 25
 PUR 278



DUAL MASS SPECTRUM
 12/28/89 20:34:00 + 5:48
 SAMPLE: 100UL C08309679 EPA#8201A CASE#18756 ON #13
 ENHANCED (5 15B 2H) *PC Salt (hhd)*
 203 BENZENE (71-43-2) NA#26

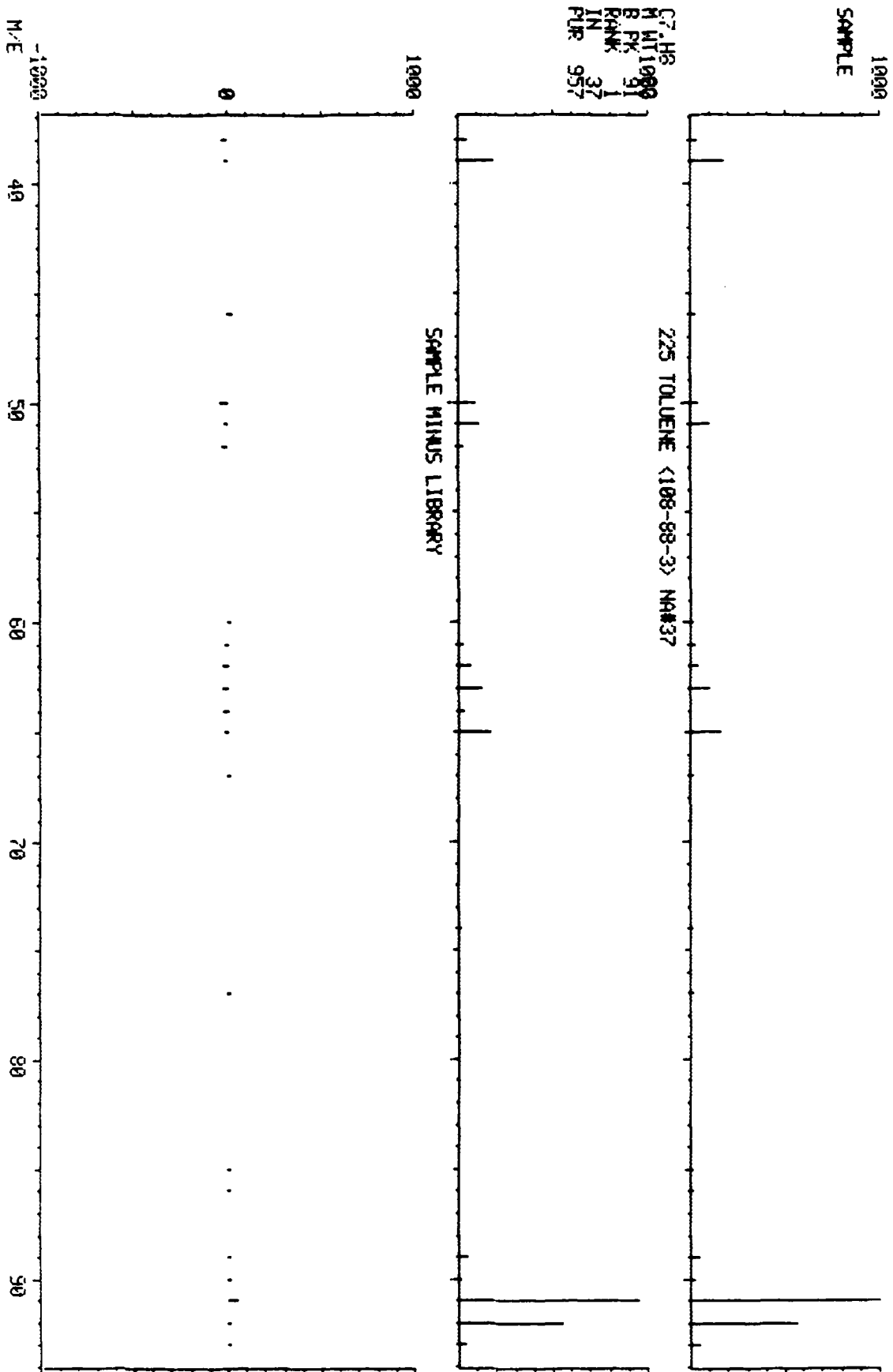
COMPUTER LABS
 DATA: C0809679B13 #464
 BASE M/E: 65/ 65
 RIC: 9375. / 21823.



LIBRARY SEARCH
12/29/89 20:34:00 + 8:52
SAMPLE: 100UL CC#309679 EPA#8201A CASE#18756 ON #13
ENHANCED (5 158 2N 0T)

COMPUCHEN LABS
RE SAN 1/1/90
DATA: C3R09679B13 # 709
BASE M/E: 91
RIC: 19039.

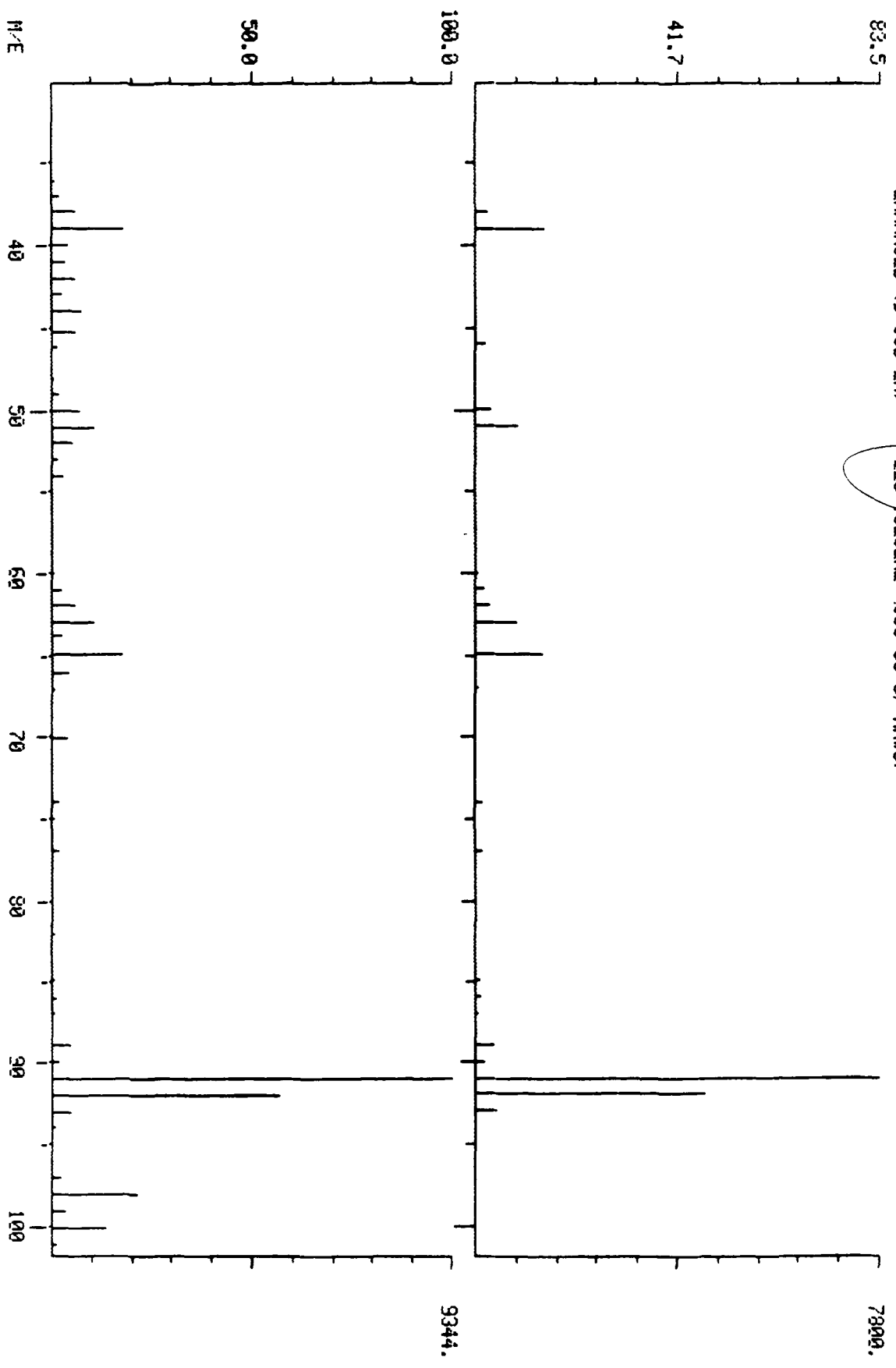
C7 H8
M W T 1000
R P X 91
R N N K 37
I N 37
P U R 957



DUPL MASS SPECTRUM
12/28/89 20:34:00 + "8:52
SAMPLE: 100UL CC#309679, EP#201A, CASE#18756 ON #13
ENHANCED (5 158 2N) 225 TOLUENE (108-88-3) NA#37

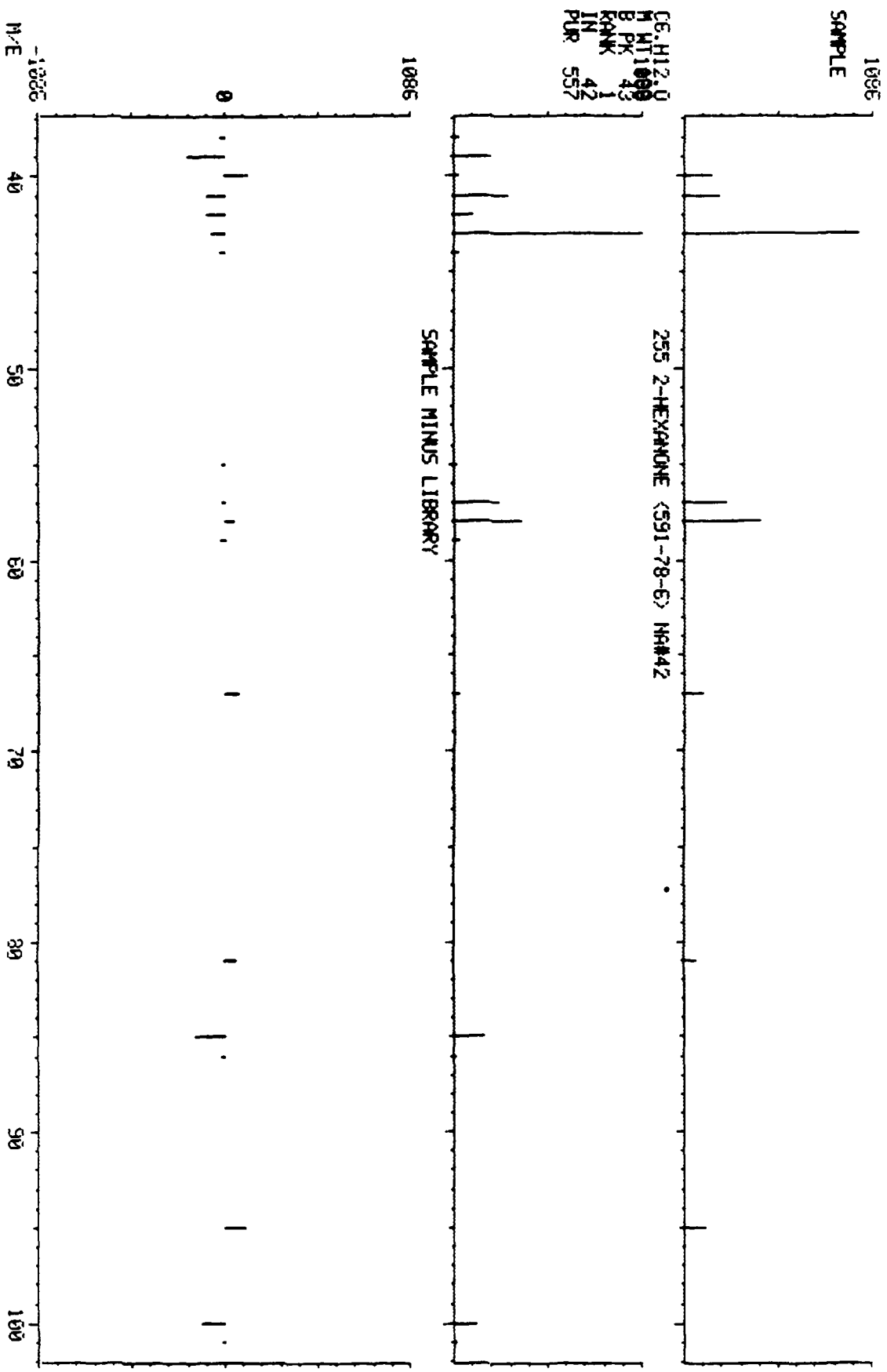
COMPUHEM LABS

DATA: C3K09679B13 #709 BASE M/E: 91 / 91
R1C: 19071. / 32639.



COMPUCHEM LIBS
 LIBRARY SEARCH
 12/28/89 20:34:00 + 10:25
 SAMPLE: 100UL CC#309679 EPA#82014, CASE#18756 ON #13
 ENHANCED (5 150 2H 01)
 DATA: C3R09679B13 # 833
 BASE M/E: 43
 RIC: 740.

CG.H12.0
 M HT 1000
 B PK 43
 KANK 1
 IN 42
 PUR 557

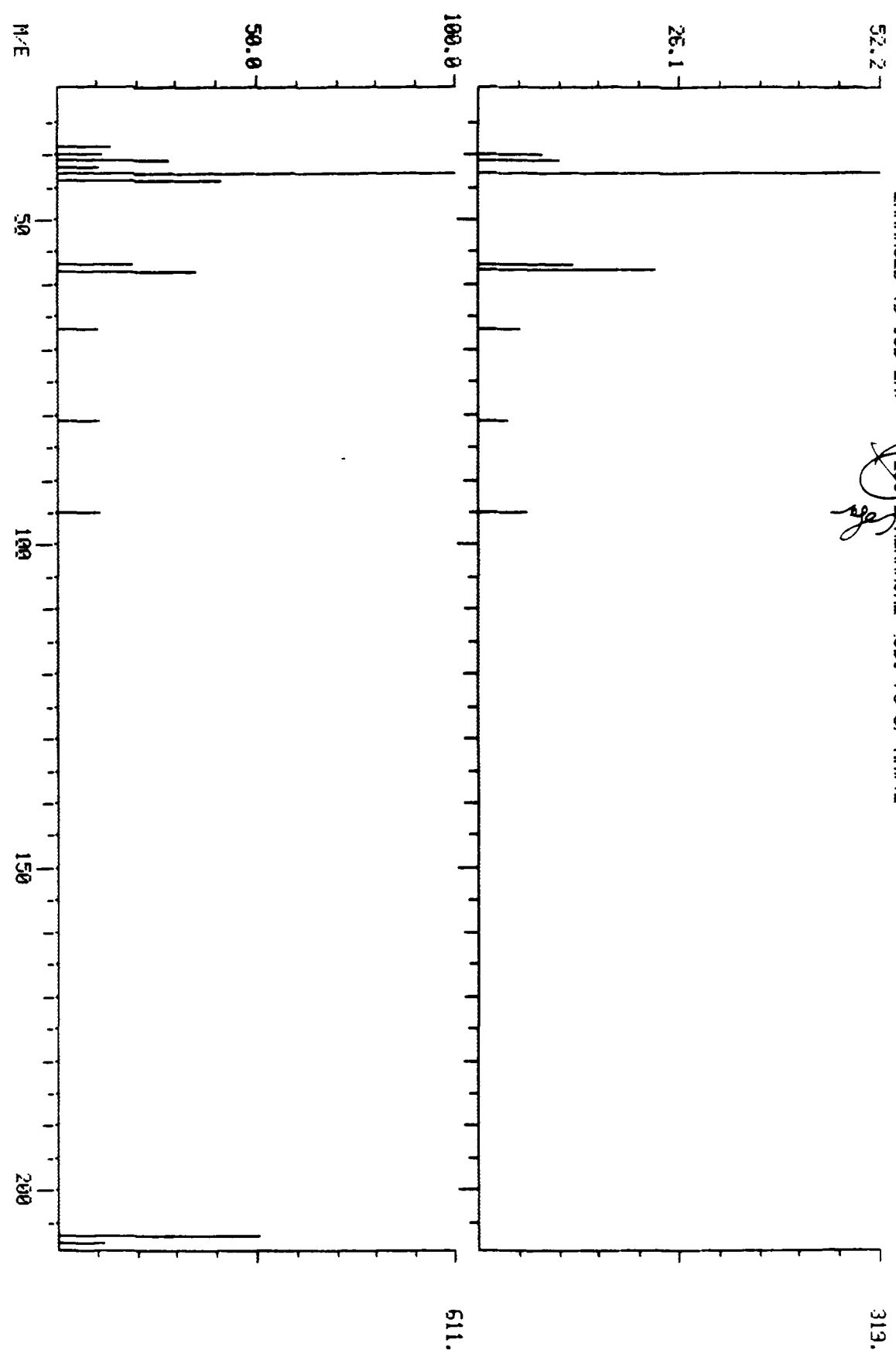


DUAL MASS SPECTRUM
12/28/89 20:34:00 + 10:25
SAMPLE: 100UL CC#309678
ENHANCED (S 158 2N)

Handwritten signature

COMFUCHEM LABS
RE 500.1440

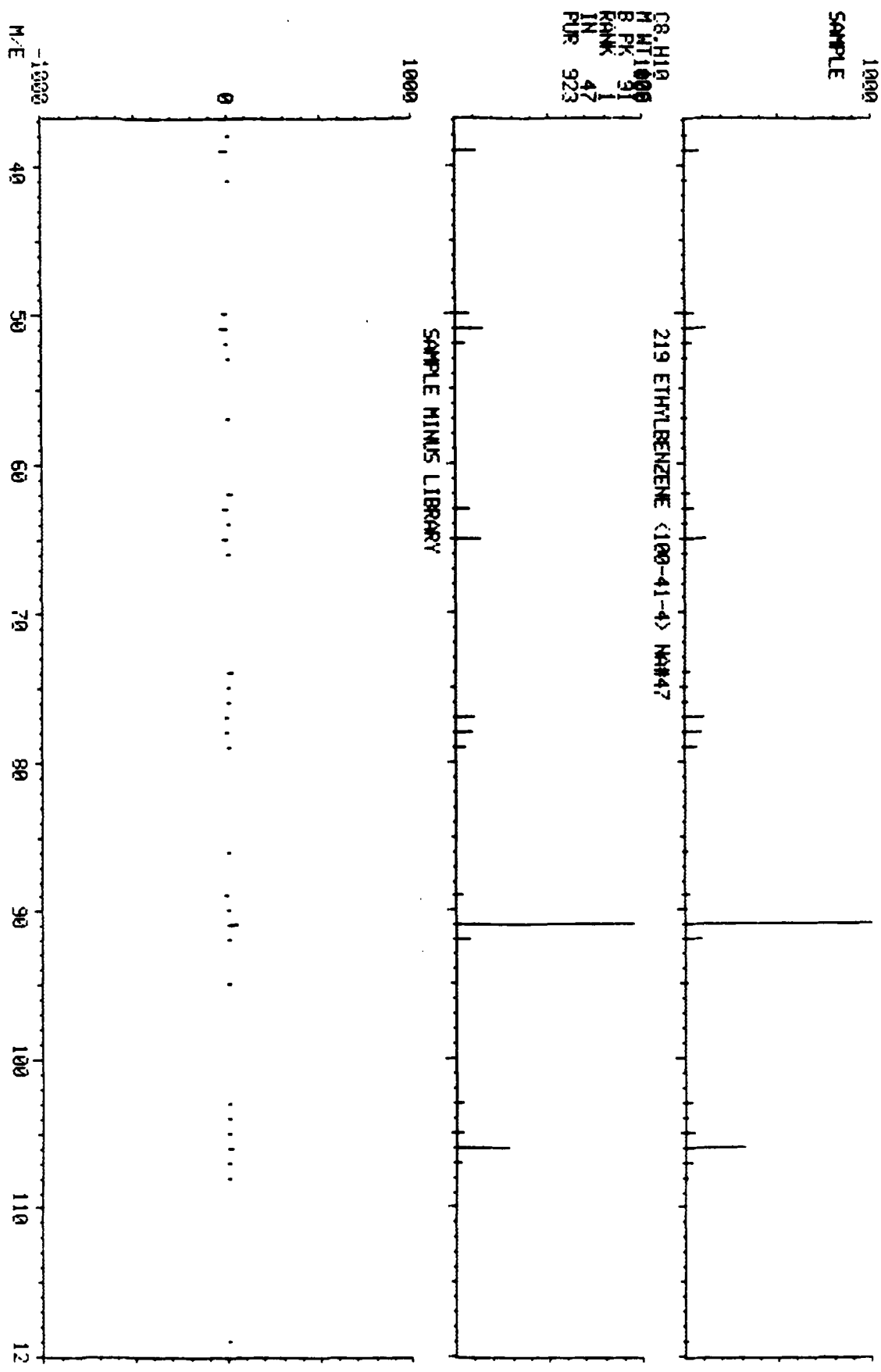
DATA: C3R09679B13 #833
BASE M/E: 43/
RIC: 740./ 2155.



COMPUchem LABS
 LIBRARY SEARCH
 12/28/89 20:34:00 + 11:37
 SAMPLE: 100UL C0309579 EPA#B201A CASE#18756 ON #13
 ENHANCED (S 15R 2N 0T)

DATA: C0309579B13 # 930
 BASE M/E: 31
 RIC: 37183.

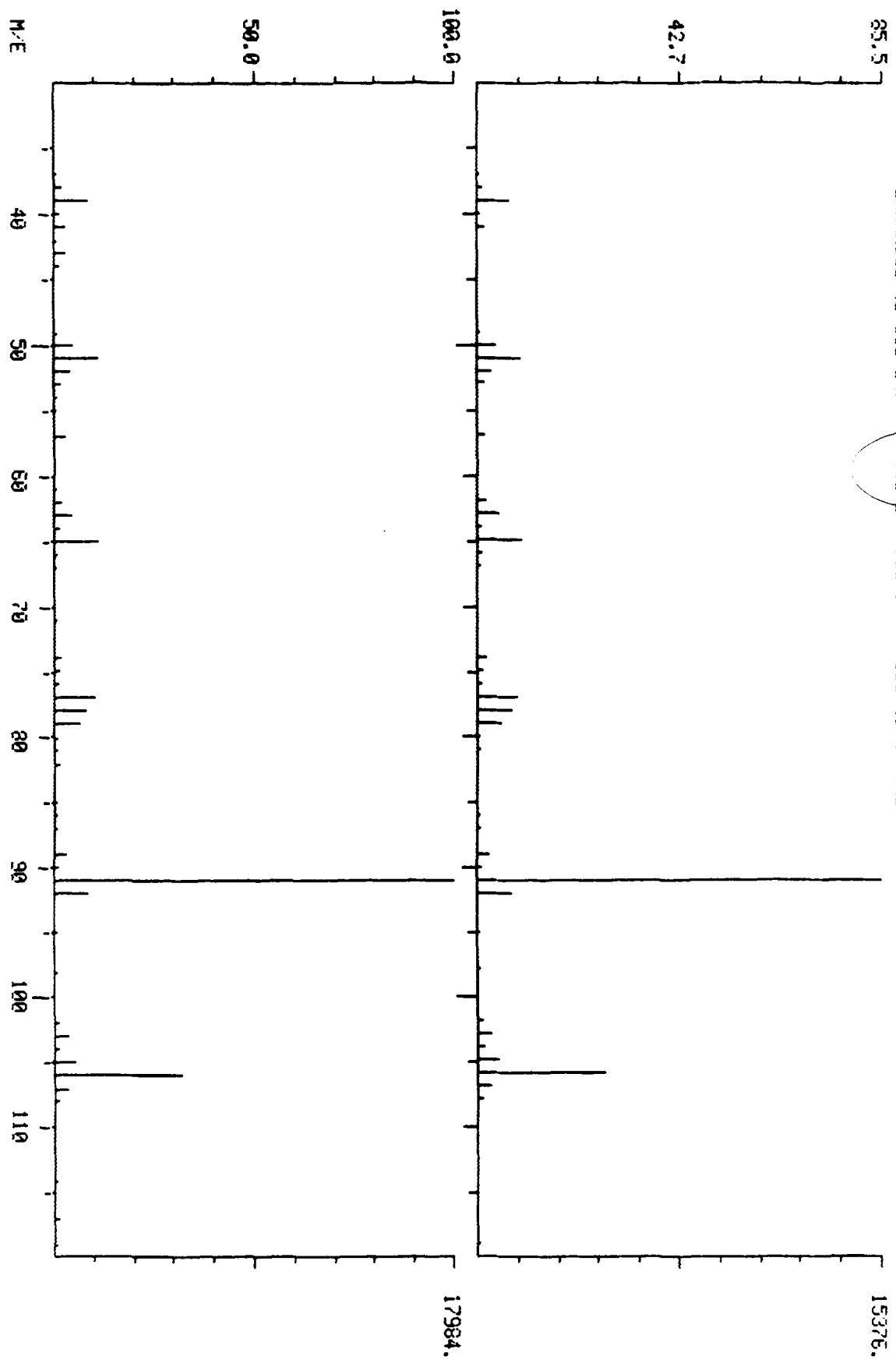
C8-H10
 M HT 1000
 B PK 31
 FRANK 47
 IN 1
 PUR 923



DUAL MASS SPECTRUM
12/28/89 20:34:00 + 11:37
SAMPLE: 100UL CC#309679 EPAN#201A CASE#18756 ON #13
ENHANCED (5 158 2M) 219 ETHYLBENZENE (100-41-4) N#47

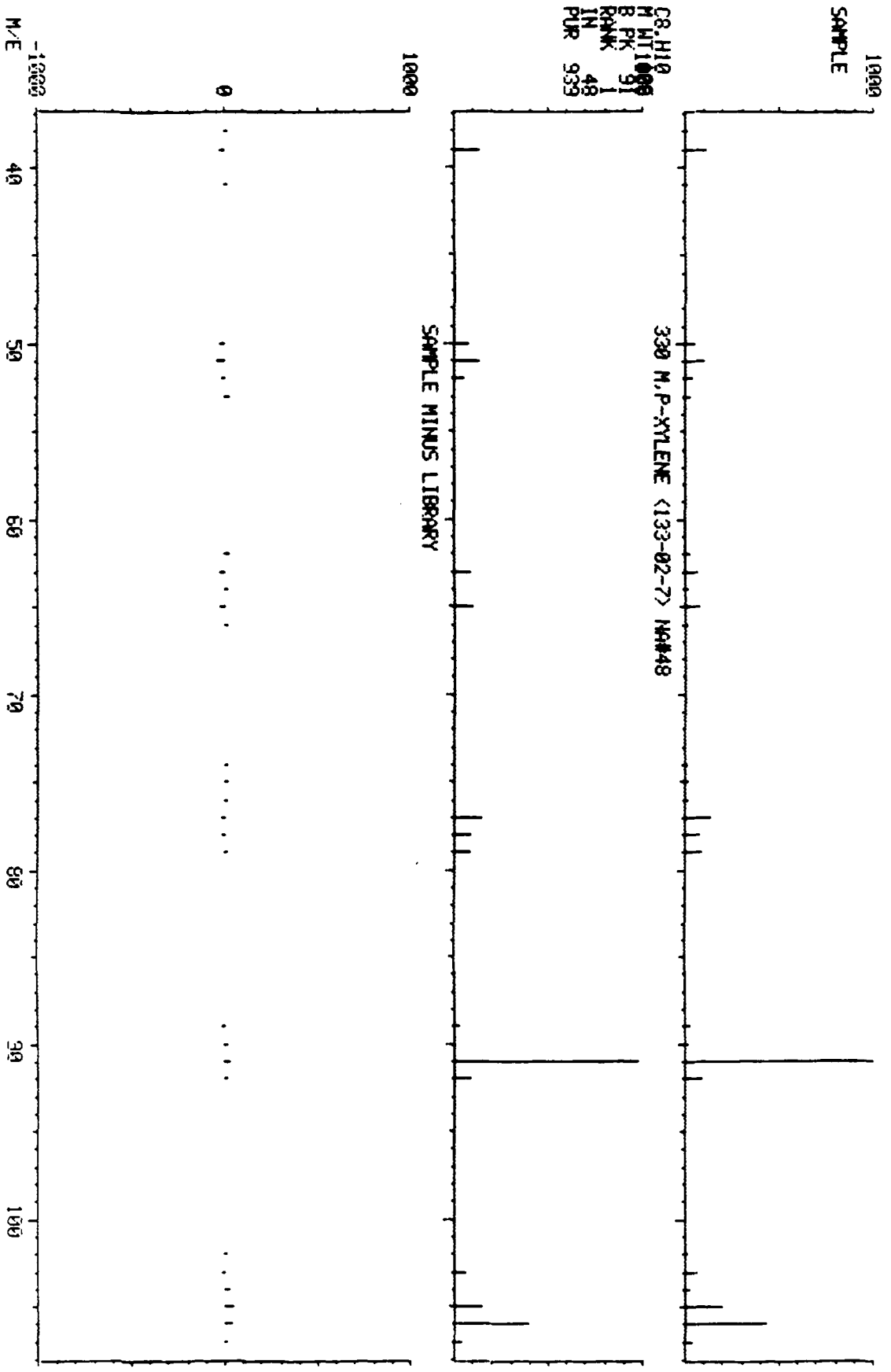
COMPUchem LABS

DATA: C3R09679B13 #330 BASE M/E: 91/ 91
RIC: 37951.7 46847.



COMPUCHEN LABS
 LIBRARY SEARCH
 12/28/89 20:34:00 + 11:52 RE sat 1/4/90
 SAMPLE: 100UL CC#309679 EPA#8201A, CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)
 DATA: C0609679B13 # 949
 BASE M/E: 91
 RIC: 86911.

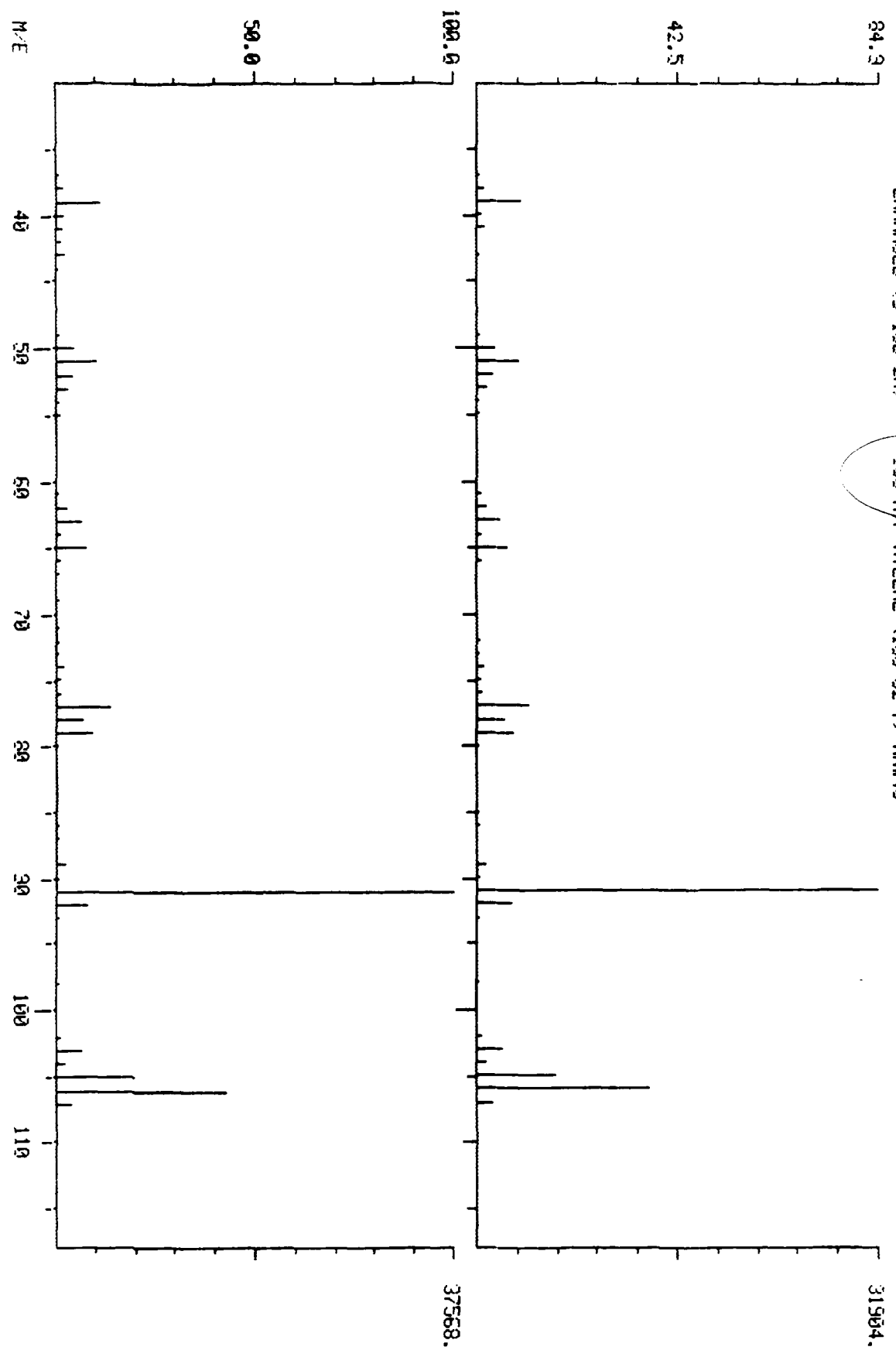
C8.H10
 M.M.I. 1000
 R.P.K. 91
 RANK 1
 IN 48
 PUR 939



DUAL MASS SPECTRUM
12/28/83 20:34:00 + 11:52
SAMPLE: 100UL CC#3909679 EPA#8201A CASE#18756 ON #13
ENHANCED (5 158 2N)

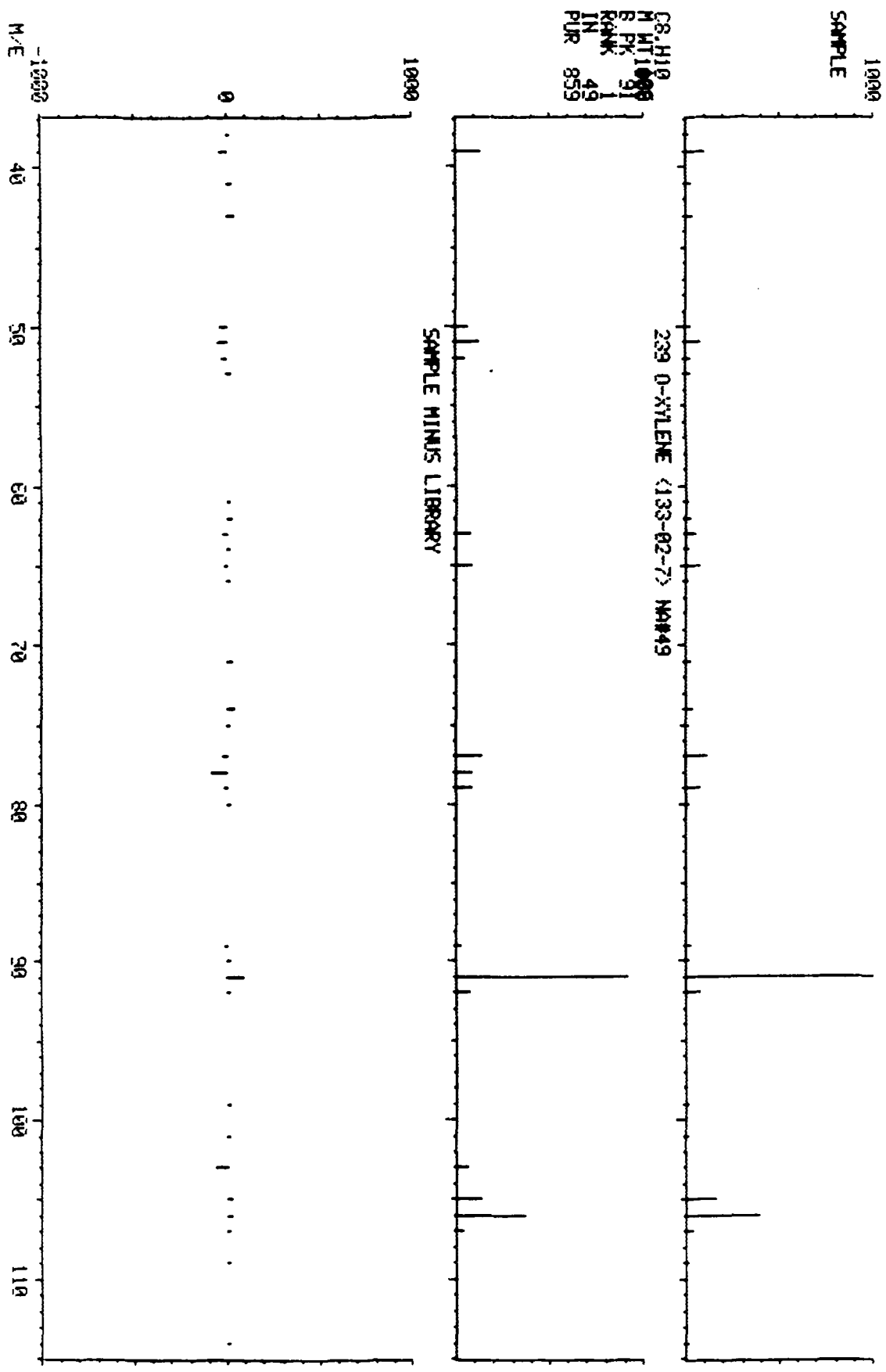
COMPUchem LABS
PE AND JJJQ
339 M/P-XYLENE (133-02-7) NA#48

DATA: 03R09679813 #349 BASE M/E: 31/ 31
RIC: 89727.7/ 109183.



COMFUCHEM LABS
 LIBRARY SEARCH
 12/28/89 20:34:00 + 12:32
 SAMPLE: 100UL CCR09579 EPA#B201Q CASE#18756 ON #13
 ENHANCED (S 15B 2H 0T)
 DATA: CCR09579B13 #1003
 BASE M/E: 31
 RIC: 51583.

CS.H10
 N.M.I. 1000
 B.P.K. 31
 K.M.K. 49
 I.N. 1
 P.U.R. 859

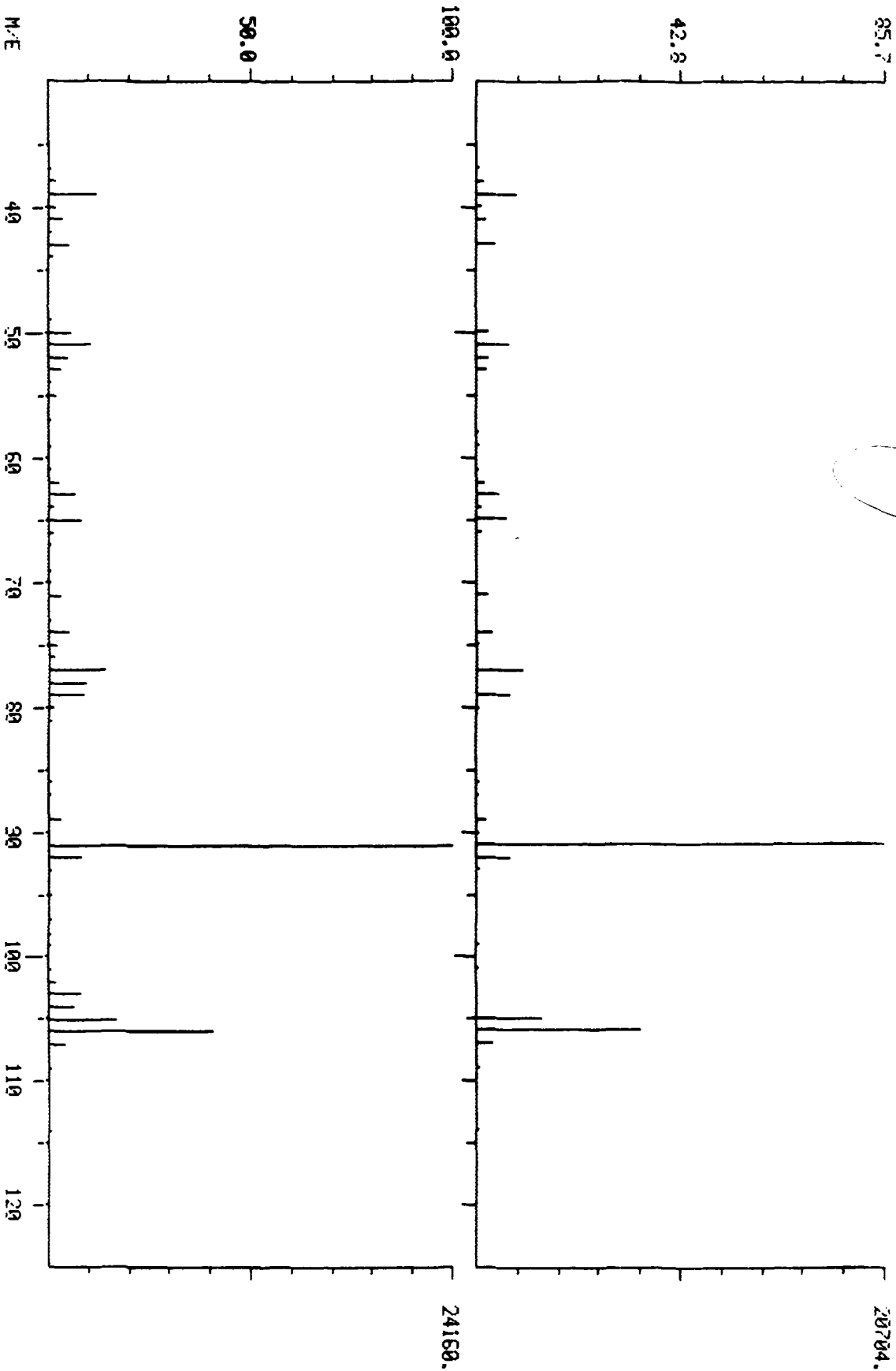


DUAL MASS SPECTRUM
12/28/89 20:34:00 + 12:32
SAMPLE: 100UL CC#309679
ENHANCED (5 158 2N)

RB LAB 11140
EPA#B201A CASE#18755 ON #13
239 D-XYLENE (133-02-7) NAW49

CONFLUENT LABS

DATA: CRR09679B13 #1003 BASE M/E: 91/ 91
RIC: 52671.7 75135.

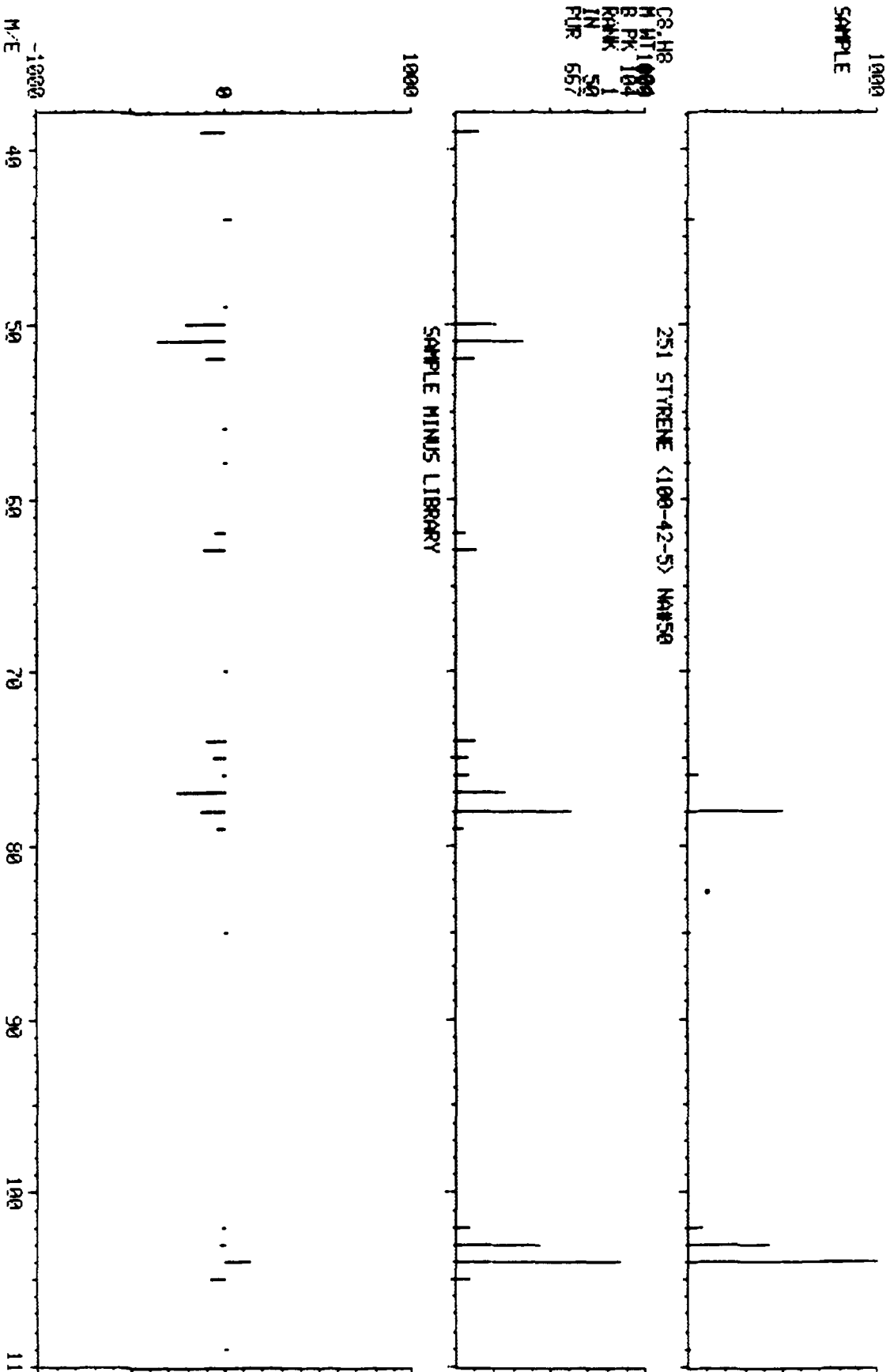


COMPUCHEM LABS
 LIBRARY SEARCH
 12/28/89 20:34:00 + 12:37
 SAMPLE: 100UL C0309579 EPA#B201A, CASE#18756 ON #13
 ENHANCED (5 158 2H 0T)

DATA: C309579B13 #1009

BASE M/E: 104
 RIC: 9967.

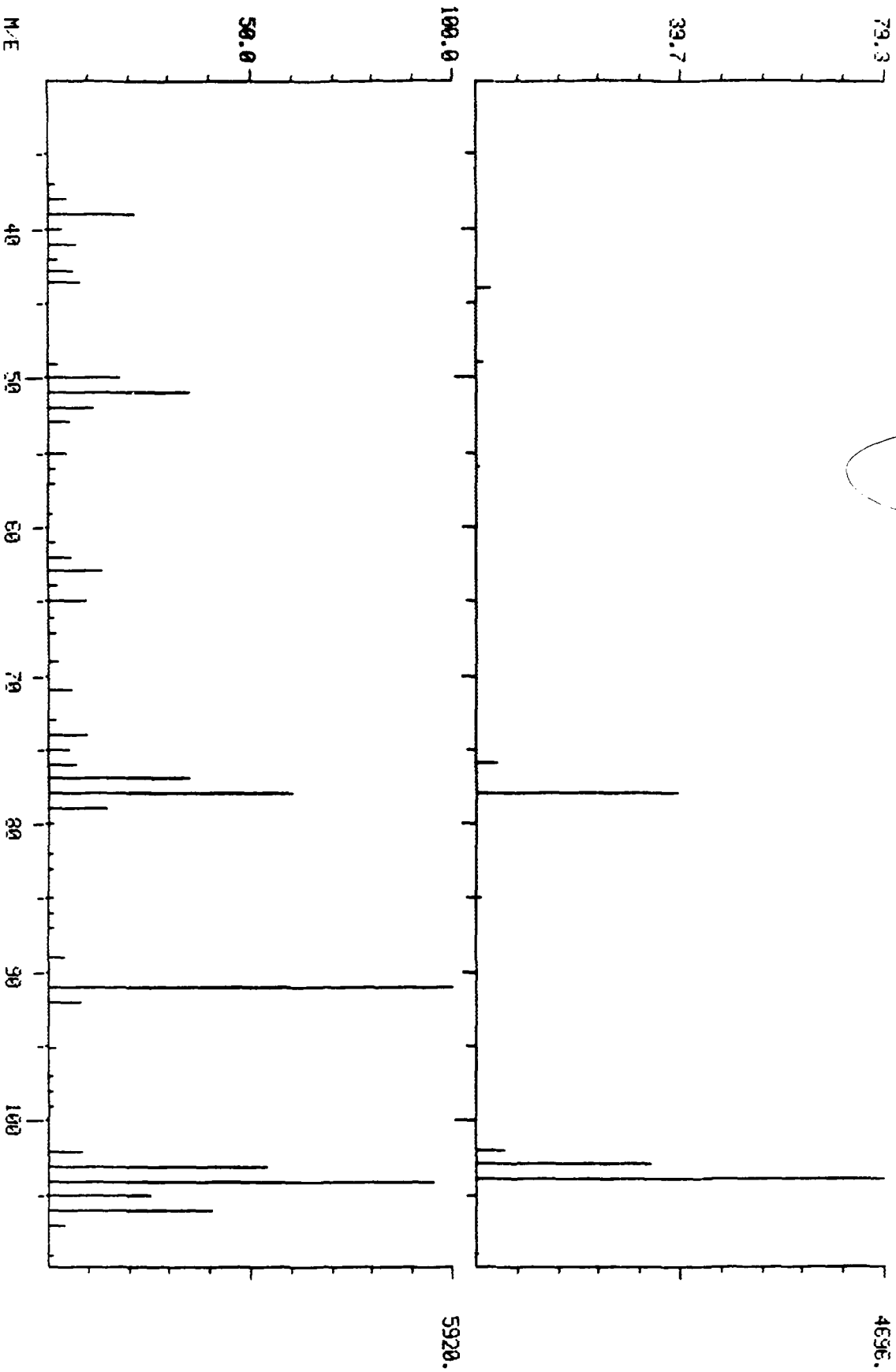
C8.H8
 H AT 1009
 R PK 104
 RANK
 IN 50
 PUR 657



DUAL MASS SPECTRUM
12/28/89 20:34:00 + 12:37
SAMPLE: 100UL CC#399679 EPA#D201A,CASE#18756 ON #13
ENHANCED (S 158 2H) 251 STYRENE <100-42-5> NA#50

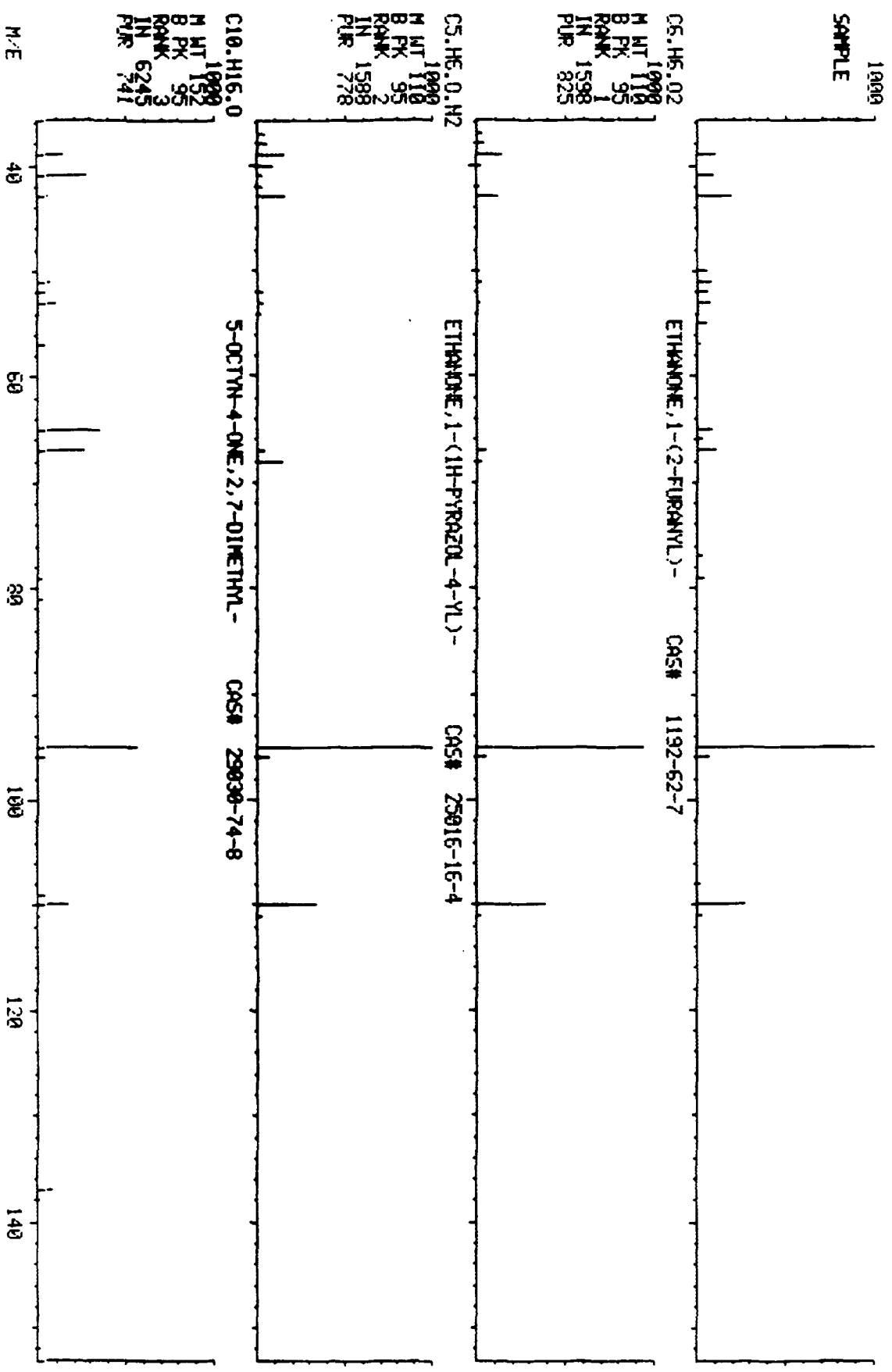
COMPLICHEM LABS
PC 2008 11/11/0

DATA: C399679813 #1009 BASE M/E: 104/ 31
RIC: 3967.7/ 39231.



LIBRARY SEARCH
 12/28/89 20:34:00 + 9:48
 SAMPLE: 100UL CC#309679 EPAB201A,CASE#18756 ON #13
 ENHANCED (S 158 2H 0T)

CONPUCHEM LABS
 DATA: C3P09679B13 # 784
 BASE M/E: 95
 RIC: 7815

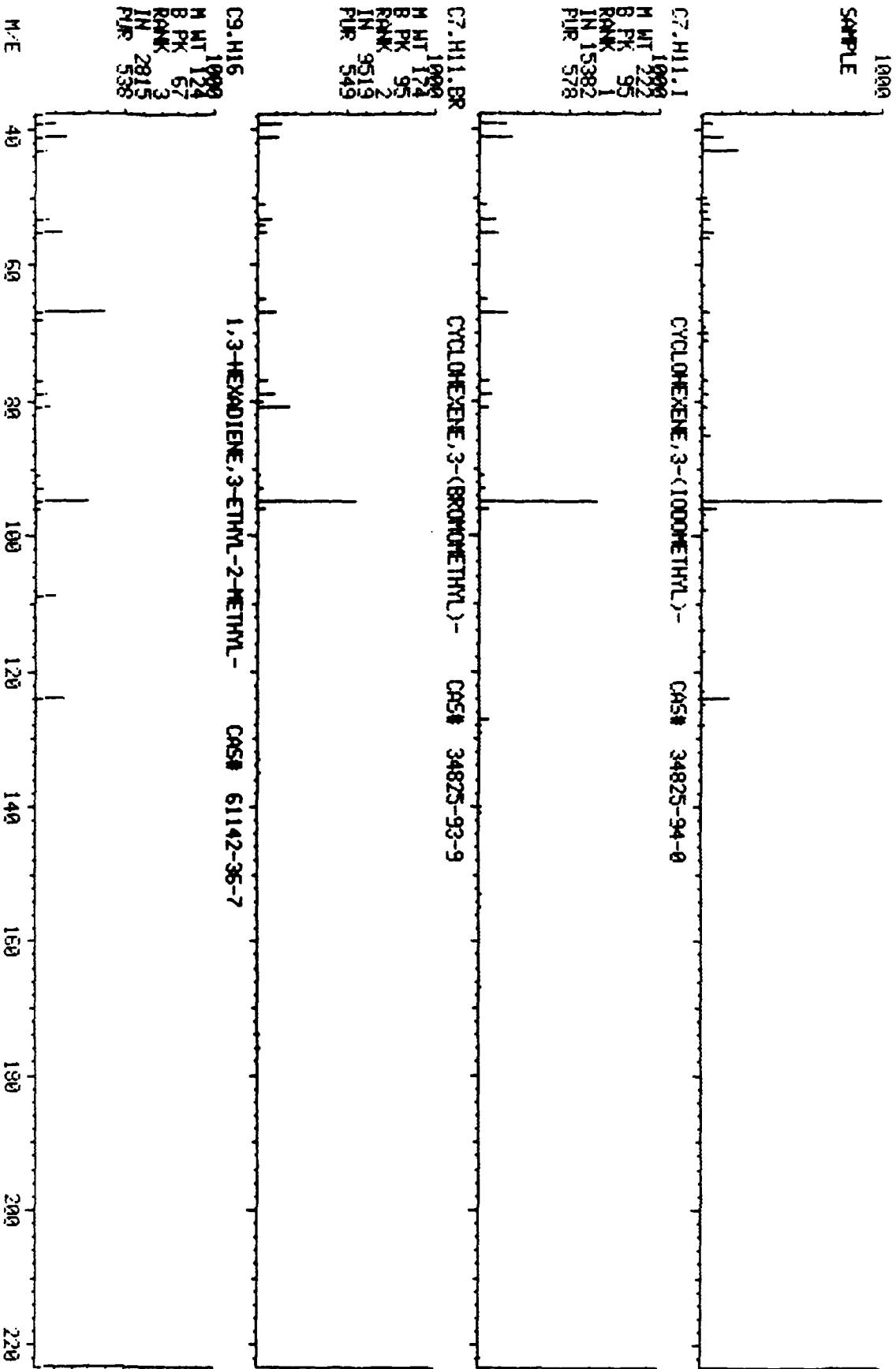


LIBRARY SEARCH
 12/28/89 20:34:00 + 12:11
 SAMPLE: 100UL CC#309679 EPA#82010, CASE#18755 ON #13
 ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

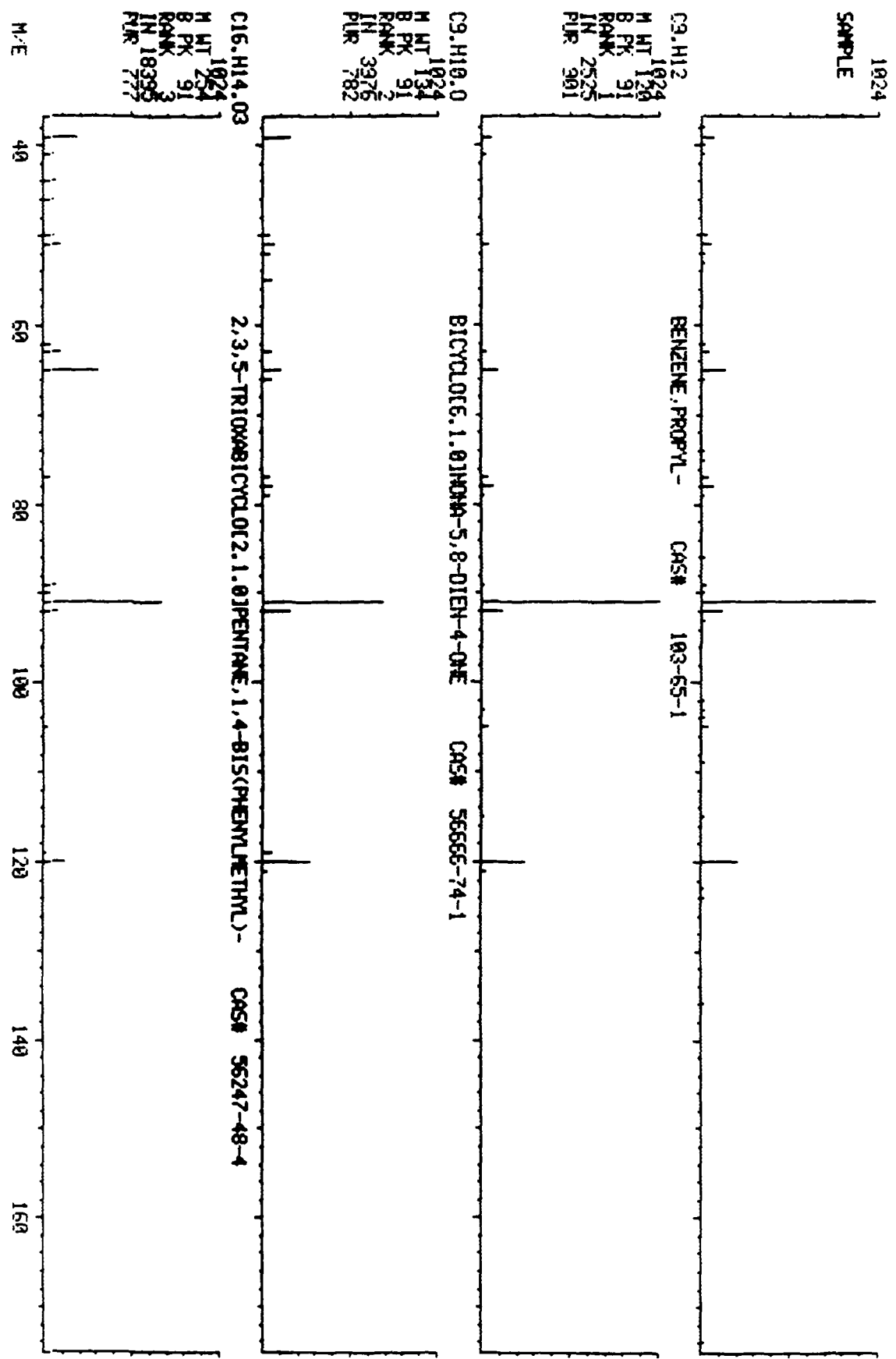
DATA: C3R09679B13 # 375

BASE M/E: 95
 RIC: 9775.



LIBRARY SEARCH
 12/28/89 20:34:00 + 14:04
 SAMPLE: 100ML CC#309579 EPABR201A CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

COMPUCHEM LABS
 DATA: C3P09579B13 #1125
 BASE M/E: 91
 RIC: 20831.

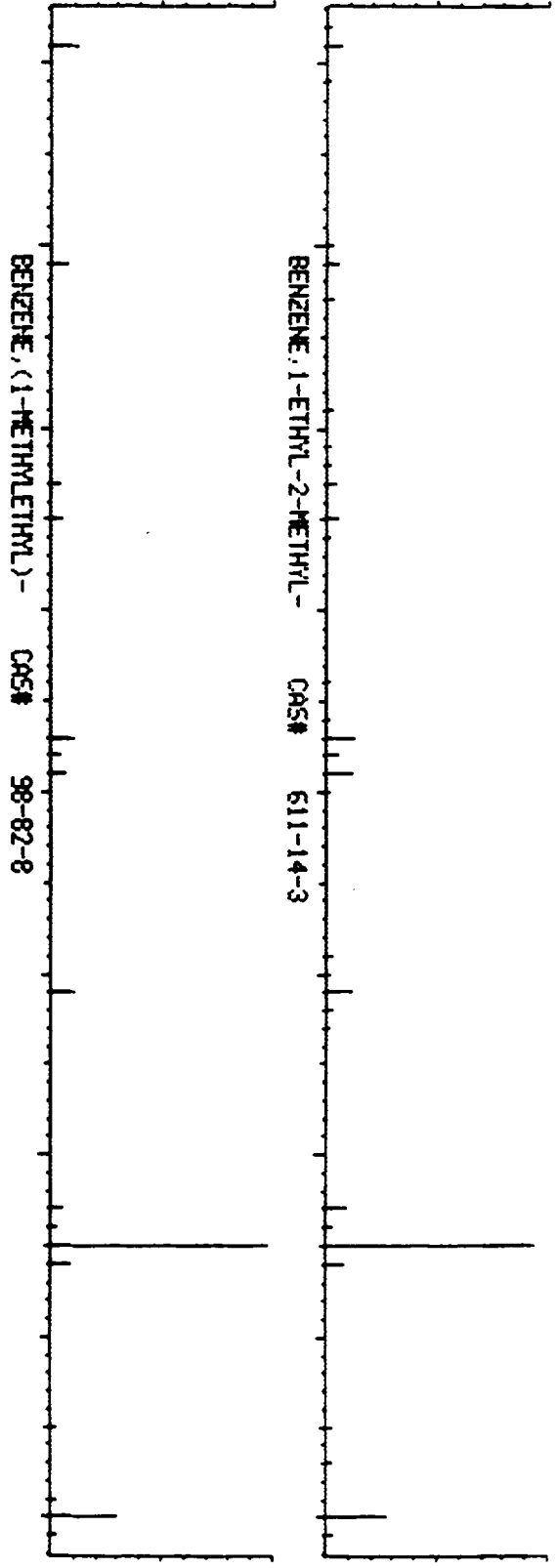


LIBRARY SEARCH
 12/28/89 20:34:00 + 14:17
 SAMPLE: 100UL CCR309579 EPA#B201A CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

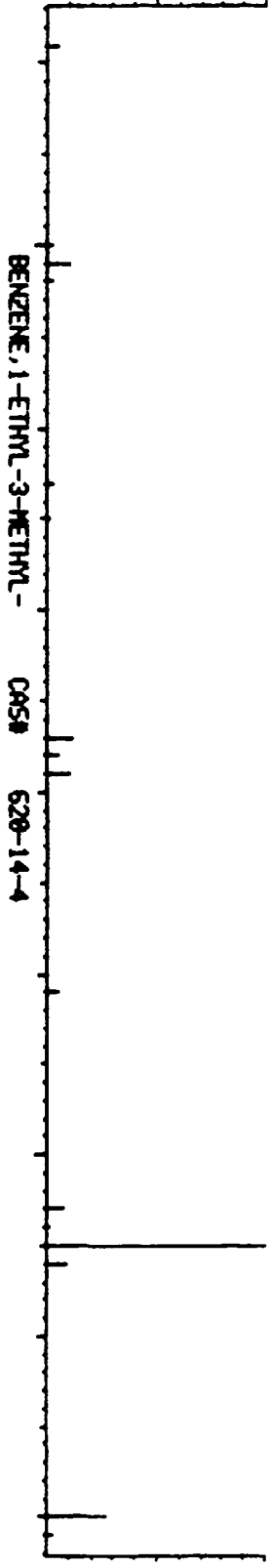
COMPUchem LABS
 RE *good* 1/1/90
 DATA: CCR309579B13 #1143
 BASE M/E: 105
 RIC: 69887.

1063
 SAMPLE

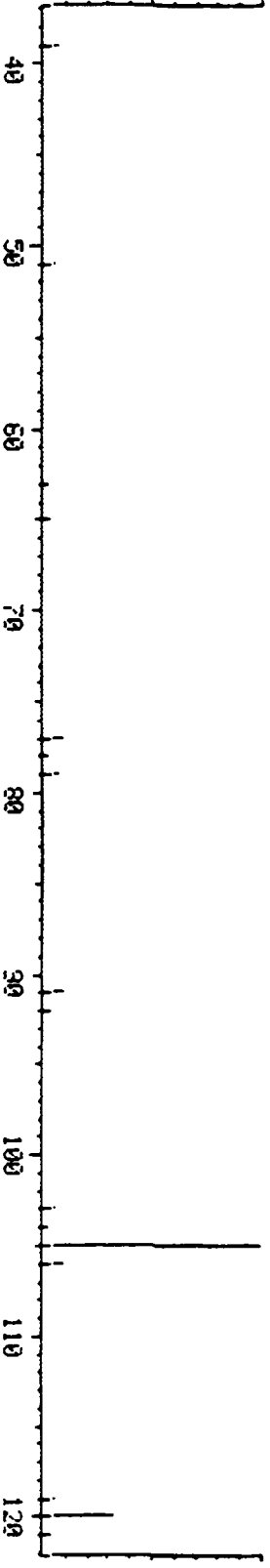
C9.H12
 1063
 M UT 120
 B PK 105
 RANK 1
 IN 2528
 PUR 300



C9.H12
 1063
 M UT 120
 B PK 105
 RANK 2
 IN 2524
 PUR 892

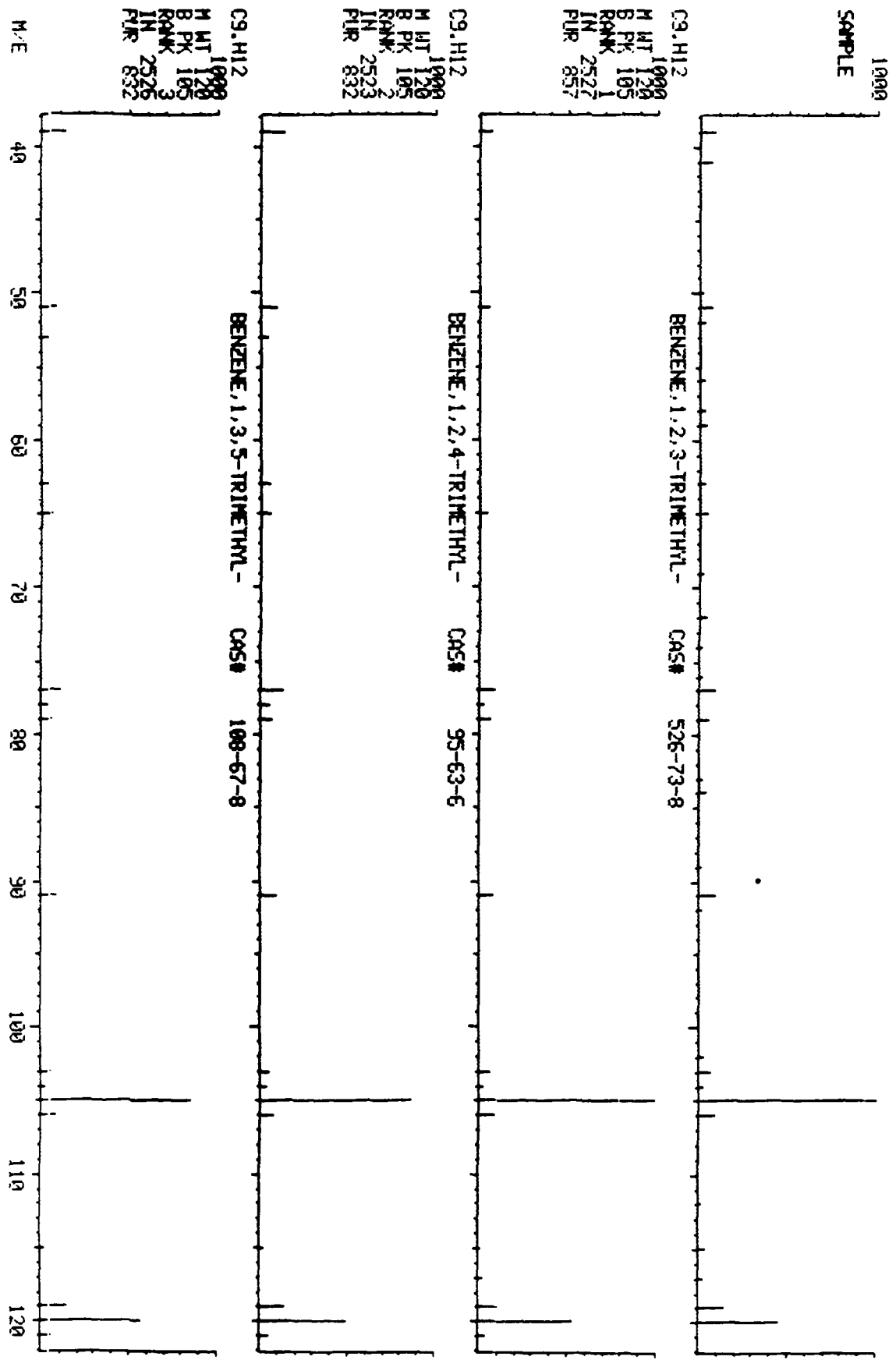


C9.H12
 1063
 M UT 120
 B PK 105
 RANK 3
 IN 2529
 PUR 883



LIBRARY SEARCH
 12/28/83 20:34:00 + 14:27
 SAMPLE: 100UL CC#309679 EPA#B201A, CASE#18756 ON #13
 ENHANCED (5 158 2N 0T)

COMPUCHEM LABS
 RE SA # 1479
 DATA: C3P09679B13 #1156
 BASE M/E: 105
 RIC: 18943.



LIBRARY SEARCH
 12/28/83 20:34:00 + 14:45 RE SA W Jhd
 SAMPLE: 100UL CC#309679 EPA#8201A CASE#18756 ON #13
 ENHANCED (S 158 2H 0T)

COMPUchem LABS

DATA: C309679B13 #1182

BASE M/E: 105
 RIC: 55935.

1075
 SAMPLE

C9.H12

M RT 1075
 B PK 105
 RANK 1
 IN 2528
 PUR 648

BENZENE, 1-ETHYL-2-METHYL-

CAS# 611-14-3

C9.H12

M RT 1075
 B PK 105
 RANK 2
 IN 2529
 PUR 639

BENZENE, 1-ETHYL-3-METHYL-

CAS# 628-14-4

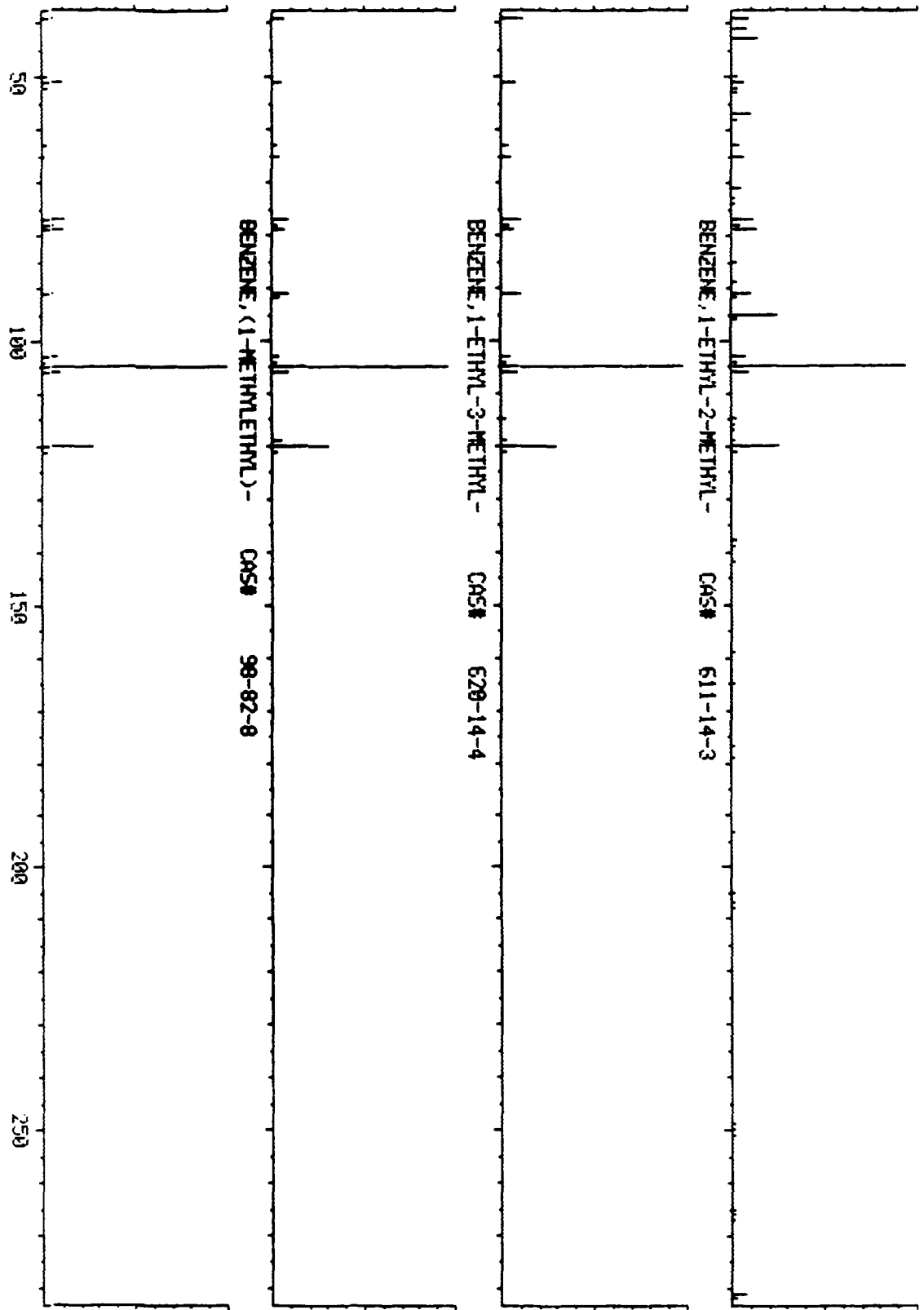
C9.H12

M RT 1075
 B PK 105
 RANK 3
 IN 2524
 PUR 635

BENZENE, (1-METHYLETHYL)-

CAS# 98-82-8

M/E

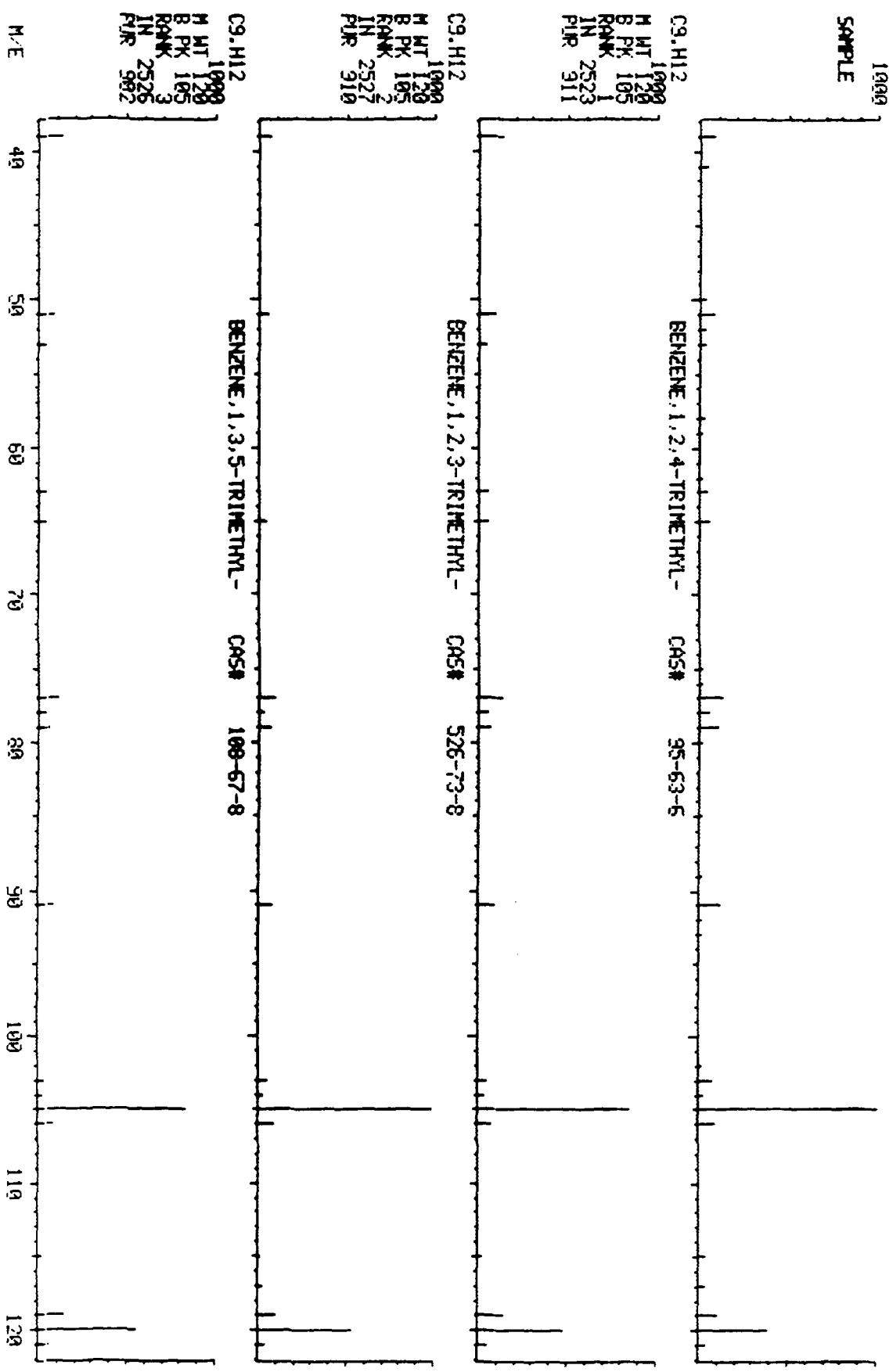


LIBRARY SEARCH
 12/28/89 20:34:08 + 15:08
 SAMPLE: 100UL CC#309579 EPA#B201A, CASE#18756 ON #13
 ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

RE *see 1/1/90* DATA: C309579B13 #1211

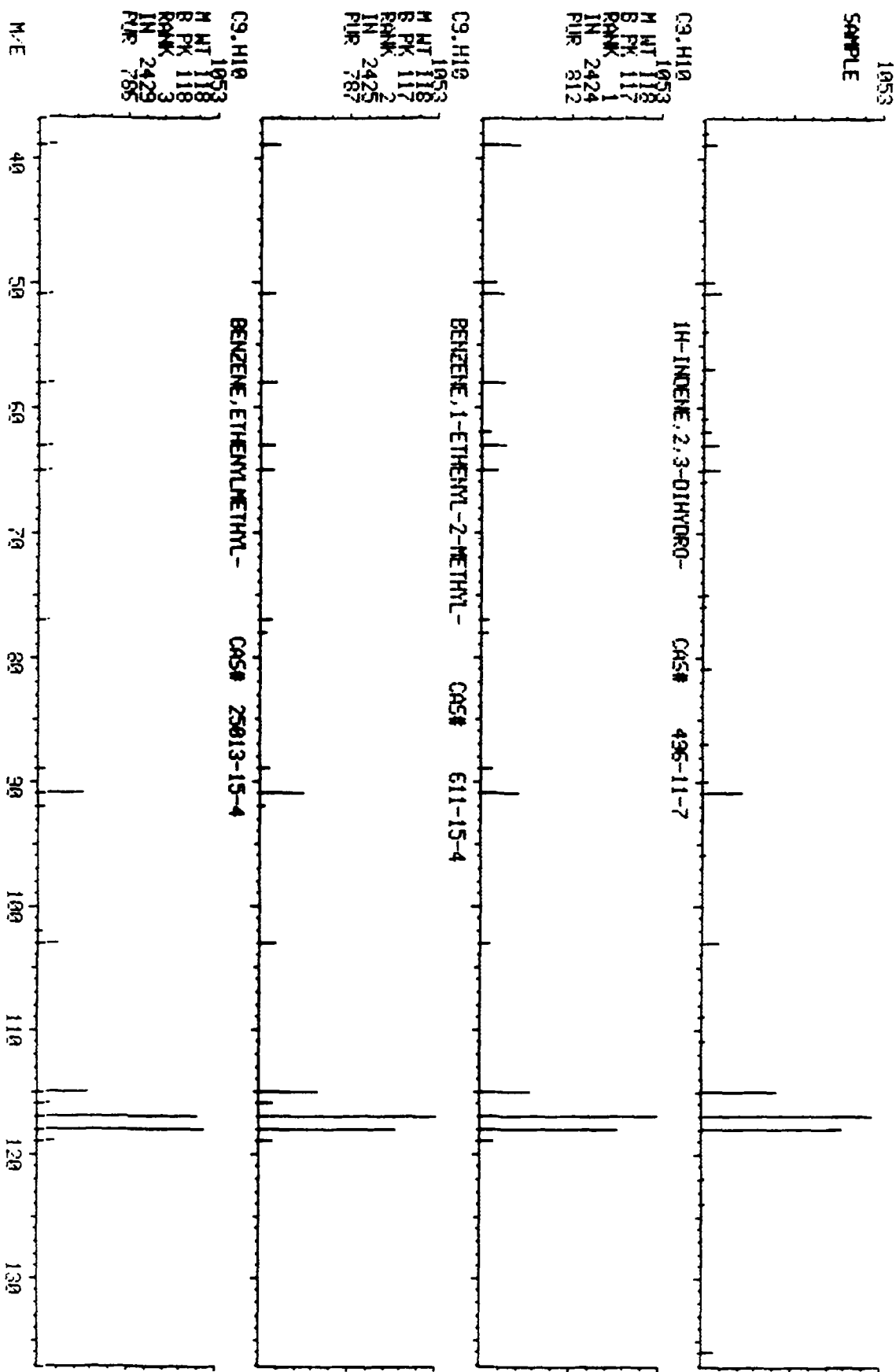
BASE M/E: 105
 RIC: 104191.



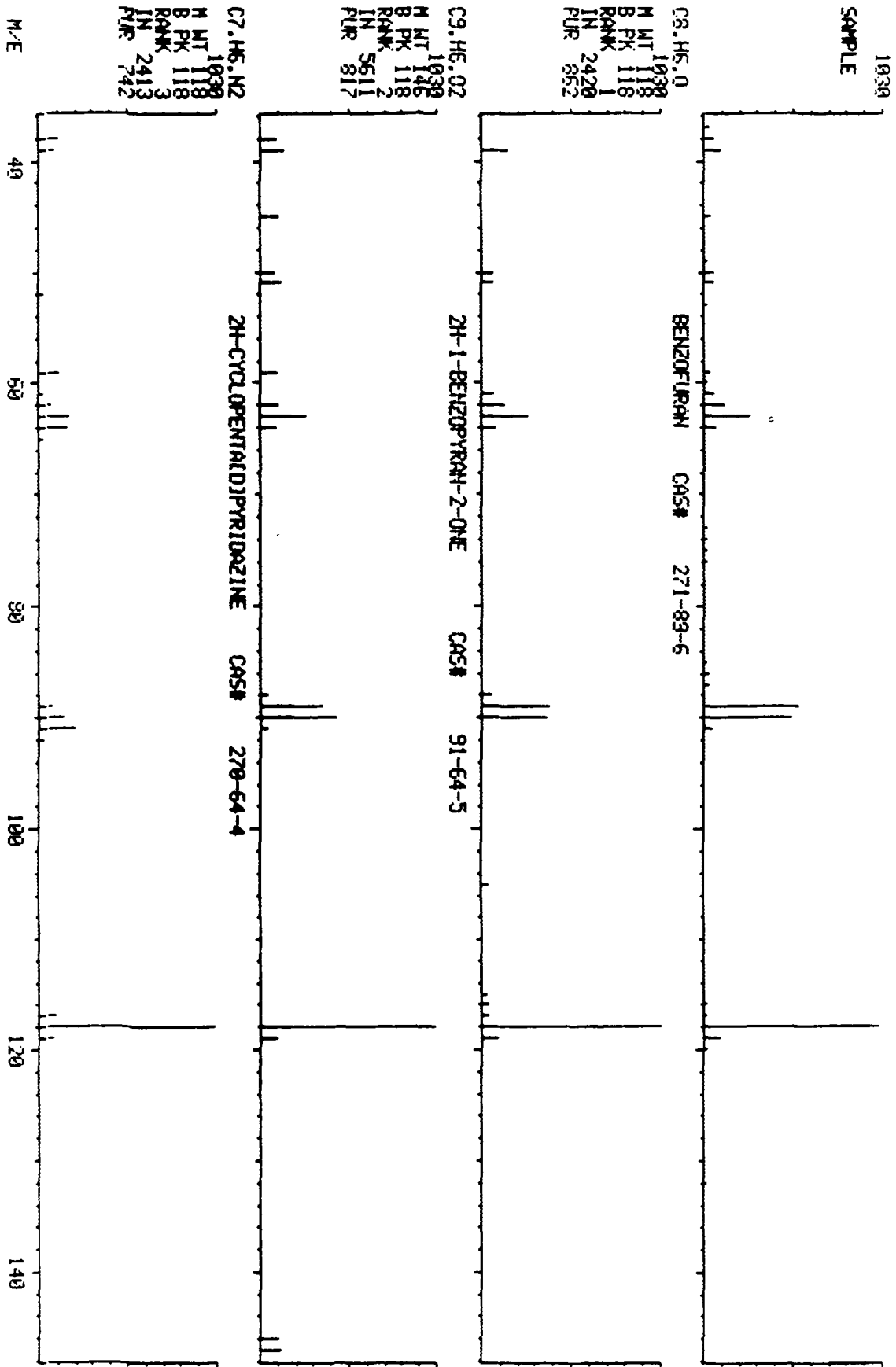
LIBRARY SEARCH
 12/28/89 20:34:00 + 15:18
 SAMPLE: 100UL CC#309679 EPA#B201A CASE#18756 ON #13
 ENHANCED (5 158 21 01)

COMPUCHEM LABS

DATA: C3R09679B13 #1224
 BASE M/E: 117
 RIC: 11743.



LIBRARY SEARCH
 12/28/89 20:34:00 + 15:33
 SAMPLE: 100UL CC#309679 EPA#B201A, CASE#18756 ON #13
 ENHANCED (5 158 2N 0T)
 COMPUTHER LABS
 DATA: CAR09679B13 #1244
 BASE M/E: 118
 RIC: 142591.



CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT. LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	399	49800	50.0		
221	50	CHLOROMETHANE				BDL	1200
231	62	VINYL CHLORIDE				BDL	1200
220	94	BROMOMETHANE				BDL	1200
209	64	CHLOROETHANE				BDL	1200
216	96	1,1-DICHLOROETHENE				BDL	620
254	76	CARBON DISULFIDE				BDL	620
252	43	ACETONE (2-PROPANONE)				BDL	1200
248	114 I	1,4-DIFLUOROBENZENE (IS)	529	183000	50.0	470	
222	84	METHYLENE CHLORIDE			3.2	4000 130	620
226	96	TRANS-1,2-DICHLOROETHENE				BDL	620
214	63	1,1-DICHLOROETHANE				BDL	620
257	43	VINYL ACETATE				BDL	1200
237	96	CIS-1,2-DICHLOROETHENE				BDL	620
253	72	2-BUTANONE				BDL	1200
211	83	CHLOROFORM				BDL	620
227	97	1,1,1-TRICHLOROETHANE				BDL	620
206	117	CARBON TETRACHLORIDE				BDL	620
203	78	BENZENE			4.6	930 5800 0	620
215	62	1,2-DICHLOROETHANE				BDL	620
270	117 I	D5-CHLOROENZENE (IS)	903	234000	50.0		
229	130	TRICHLOROETHENE				BDL	620
217	63	1,2-DICHLOROPROPANE				BDL	620
212	83	BROMODICHLOROMETHANE				BDL	620
218	75	CIS-1,3-DICHLOROPROPENE				BDL	620
256	43	4-METHYL-2-PENTANONE				BDL	1200
225	92	TOLUENE			30.3	1200 3800 0	620
250	75	TRANS-1,3-DICHLOROPROPENE				BDL	620
228	97	1,1,2-TRICHLOROETHANE				BDL	620
224	164	TETRACHLOROETHENE				BDL	620
255	43	2-HEXANONE			4.6	4.6 630 320 5800 0 1200	1200
208	129	DIBROMOCHLOROMETHANE				BDL	620
207	112	CHLOROBENZENE				BDL	620
219	106	ETHYLBENZENE			22.4	310 2800 0	620
330	106	M, P-XYLENE			40.3	5600 3000 0	620
239	106	O-XYLENE			29.0	4000 2600 0	620
251	104	STYRENE			9.5	1300 1200 0	620
205	173	BROMOFORM				BDL	620
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	620
258	65 S	D4-1,2-DICHLOROETHANE NA#57			55.5	111. %	
247	95 S	BROMOFLUOROBENZENE			46.4	93. %	
233	98 S	D8-TOLUENE NA#59			45.6	91. %	
289	106	XYLENES (TOTAL)			69.3	8700 9600 0	620


CORRECTED/REVIEWED BY

(GC/MS DATA REVIEWER)

DATE

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
299	96	1,2-DICHLOROETHENE (TOTAL)				BDL	620
CHECKSUMS:							
		3979.	1831	466800.		510.7	26955.

CORRECTED/REVIEWED BY



(GC/MS DATA REVIEWER)

DATE

11/3/90

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P	F
40	258	D4-1,2-DICHLOROETHANE NA#57	55.5	50.0	111.	70-121	X	
41	247	BROMOFLUOROBENZENE	46.4	50.0	93.	74-121	X	
42	233	D8-TOLUENE NA#59	45.6	50.0	91.	81-117	X	

* ADVISORY SURROGATE ONLY
 ++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

CORRECTION FACTOR CALCULATION:

$$\frac{4.0 \text{ g}}{\text{WET WEIGHT OF SAMPLE (g)}} \times \text{LIBRARY ADJUSTMENT} \times \text{DRY WEIGHT FACTOR} \times \frac{10000 \text{ UL}}{\text{UL USED}} =$$

$$\frac{4.0 \text{ g}}{4.00 \text{ (g)}} \times 1.25 \times \frac{1.10}{1.00} \times \frac{10000 \text{ UL}}{100. \text{ UL}} = 125.000$$

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:

$$\frac{100. \text{ UL}}{\text{UL USED}} = 1.0$$

VERSION 8

CORRECTED/REVIEWED BY

(GC/MS DATA REVIEWER)

DATE

[Handwritten signature]
 1/29/90

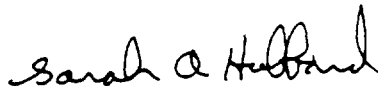
LABORATORY NOTICE
COMPUCHEM ID # 309679
CLIENT ID # B201A
CASE # 18756

The volatile fraction of this sample was prepared and analyzed as both a Low Level Solid and a Medium Level Solid. The initial analysis of 5 g sample from the Low Level preparation indicated that a Medium Level analysis was necessary in order to bring target compound concentrations into the instruments' analytical range. The analysis of 100 ul of the Medium Level extract proved to be an acceptable dilution. However, the following table of results shows that the concentrations for these analytes did not compare well.

Target Compound	Reported Concentration, ug/Kg	
	5g Low Level	100 ul Medium Level Extract
toluene	720 *	4200
ethylbenzene	420	3100
m,p-xylene	720 *	5600
o-xylene	570	4000
styrene	210	1300

*estimated concentration of saturated compound

We have concluded that these differences are due to a combination of sample inhomogeneity and the difference between the Low Level purge and trap and the Medium Level extraction procedure. In addition, the surrogate D8-toluene failed recovery acceptance criteria in the Low Level analysis due to the presence of a large concentration of the native compound. Secondary ion calculations did not improve the recovery. The data is being reported with reference to this qualifier.



Sarah A. Hubbard
Senior Volatile Data Specialist
January 3, 1990

QA Approval #1013
Linda Fowler
Employee #820
Sr. Quality Assurance Specialist
January 4, 1990

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: 309686

Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009686C19

Level: (low/med) LOW Date Received: 12/20/89

% Moisture: not dec. 12 Date Analyzed: 12/27/89

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	57	U
74-83-9	-----Bromomethane	57	U
75-01-4	-----Vinyl Chloride	57	U
75-00-3	-----Chloroethane	57	U
75-09-2	-----Methylene Chloride	69	B
67-64-1	-----Acetone	370	B
75-15-0	-----Carbon Disulfide	28	U
75-35-4	-----1,1-Dichloroethene	28	U
75-34-3	-----1,1-Dichloroethane	28	U
540-59-0	-----1,2-Dichloroethene (total)	28	U
67-66-3	-----Chloroform	28	U
107-06-2	-----1,2-Dichloroethane	28	U
78-93-3	-----2-Butanone	260	U
71-55-6	-----1,1,1-Trichloroethane	28	U
56-23-5	-----Carbon Tetrachloride	28	U
108-05-4	-----Vinyl Acetate	57	U
75-27-4	-----Bromodichloromethane	28	U
78-87-5	-----1,2-Dichloropropane	28	U
10061-01-5	-----cis-1,3-Dichloropropene	28	U
79-01-6	-----Trichloroethene	28	U
124-48-1	-----Dibromochloromethane	28	U
79-00-5	-----1,1,2-Trichloroethane	28	U
71-43-2	-----Benzene	140	U
10061-02-6	-----Trans-1,3-Dichloropropene	28	U
75-25-2	-----Bromoform	28	U
108-10-1	-----4-Methyl-2-Pentanone	71	U
591-78-6	-----2-Hexanone	1600	BE
127-18-4	-----Tetrachloroethene	28	U
79-34-5	-----1,1,2,2-Tetrachloroethane	28	U
108-88-3	-----Toluene	1600	E
108-90-7	-----Chlorobenzene	28	U
100-41-4	-----Ethylbenzene	980	U
100-42-5	-----Styrene	540	U
1330-20-7	-----Total Xylenes	3300	E

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009686C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/27/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10 CONCENTRATION UNITS;
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	11.04	500	J
2.	METHYLESTERPENTENOIC ACID +	13.45	1400	J
3.	UNKNOWN	14.09	1100	J
4. 110-43-0	2-HEPTANONE	14.27	1400	J
5.	UNKNOWN	14.42	910	J
6.	UNKNOWN	14.55	1100	J
7. 100-66-3	BENZENE, METHOXY-	14.72	2300	J
8.	UNKNOWN	15.35	500	J
9.	ETHYLMETHYLBENZENE + UNKNOWN	15.60	5400	J
10. 611-14-3	BENZENE, 1-ETHYL-2-METHYL-	16.07	1500	J

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009686C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/27/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG	Q
74-87-3	Chloromethane	57	U
74-83-9	Bromomethane	57	U
75-01-4	Vinyl Chloride	57	U
75-00-3	Chloroethane	57	U
75-09-2	Methylene Chloride	69	B
67-64-1	Acetone	370	B
75-15-0	Carbon Disulfide	28	U
75-35-4	1,1-Dichloroethene	28	U
75-34-3	1,1-Dichloroethane	28	U
540-59-0	1,2-Dichloroethene (total)	28	U
67-66-3	Chloroform	28	U
107-06-2	1,2-Dichloroethane	28	U
78-93-3	2-Butanone	260	
71-55-6	1,1,1-Trichloroethane	28	U
56-23-5	Carbon Tetrachloride	28	U
108-05-4	Vinyl Acetate	57	U
75-27-4	Bromodichloromethane	28	U
78-87-5	1,2-Dichloropropane	28	U
10061-01-5	cis-1,3-Dichloropropene	28	U
79-01-6	Trichloroethene	28	U
124-48-1	Dibromochloromethane	28	U
79-00-5	1,1,2-Trichloroethane	28	U
71-43-2	Benzene	140	
10061-02-6	Trans-1,3-Dichloropropene	28	U
75-25-2	Bromoform	28	U
108-10-1	4-Methyl-2-Pentanone	71	
591-78-6	2-Hexanone	1600	BE
127-18-4	Tetrachloroethene	28	U
79-34-5	1,1,2,2-Tetrachloroethane	28	U
108-88-3	Toluene	1600	E
108-90-7	Chlorobenzene	28	U
100-41-4	Ethylbenzene	980	
100-42-5	Styrene	540	
1330-20-7	Total Xylenes	3300	E

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B201B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009686C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/27/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	11.04	500	J
2.	METHYLESTERPENTENOIC ACID +	13.45	1400	J
3.	UNKNOWN	14.09	1100	J
4. 110-43-0	2-HEPTANONE	14.27	1400	J
5.	UNKNOWN	14.42	910	J
6.	UNKNOWN	14.55	1100	J
7. 100-66-3	BENZENE, METHOXY-	14.72	2300	J
8.	UNKNOWN	15.35	500	J
9.	ETHYLMETHYLBENZENE + UNKNOWN	15.60	5400	J
10. 611-14-3	BENZENE, 1-ETHYL-2-METHYL-	16.07	1500	J

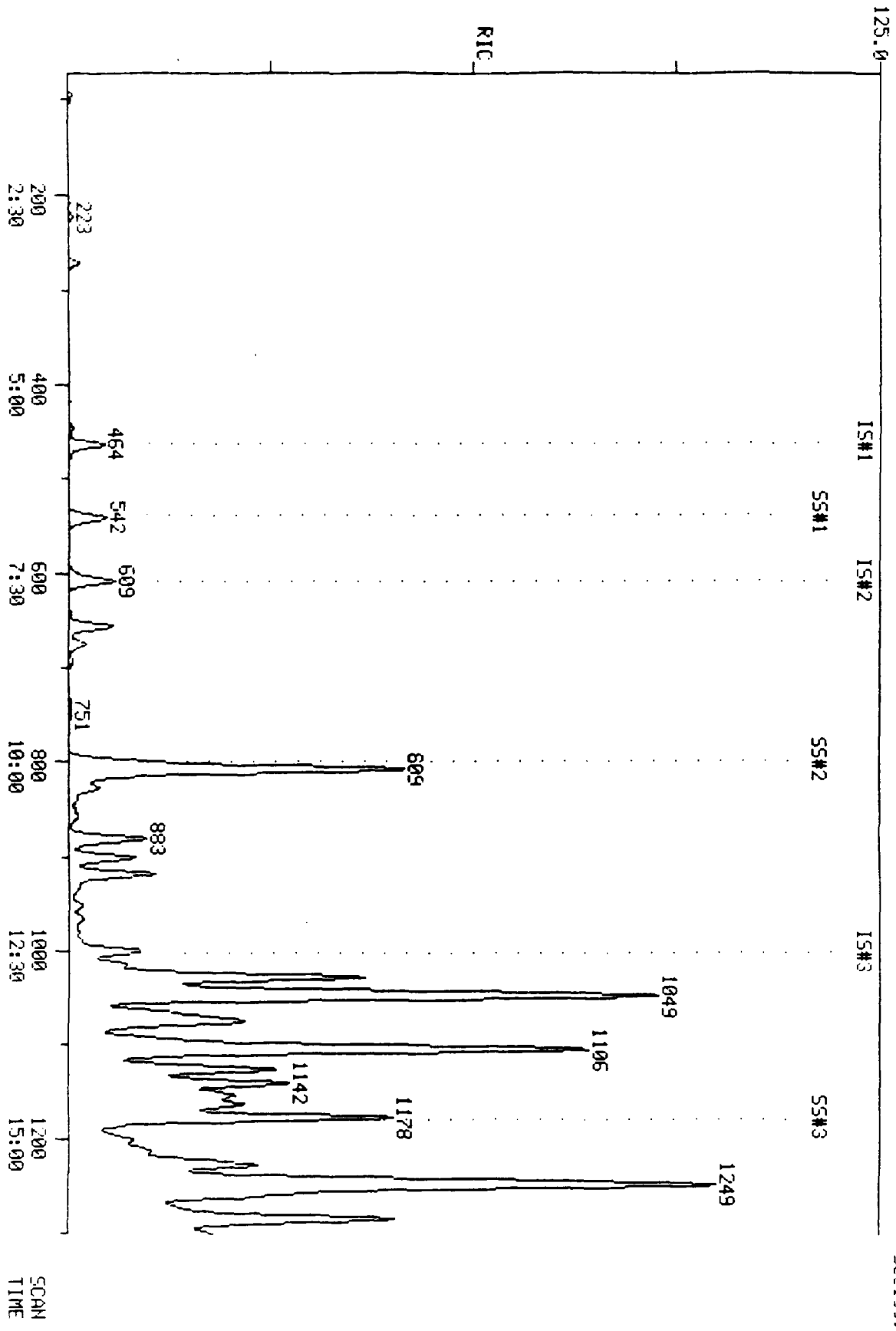
FORM I VOA-TIC

1/87 Rev.

RIC
 12/27/89 5:39:00
 SAMPLE: 1G LC#309686 CASE#18756.7 EPA#6201B ON#19
 COND5.:

COMPUchem LABS
 COMPUchem DATA: GR009686C19 SCANS 73 TO 1300

956150.



QUANTITATION REPORT FILE: GRC09686C19
 DATA: GR009686C19.TI
 12/27/89 5:39:00.
 SAMPLE: 1G CC#309686 CASE#18756 7 EPA#B201B DN#19
 CONDS.:
 SUBMITTED BY 19 ANALYST: 1422

AMOUNT=AREA * REF.AMNT/(REF.AREA)* RESP.FACT)
 RESP. FAC. FROM LIBRARY ENTRY ✓

- | NO | NAME |
|----|--|
| 1 | *234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1 |
| 2 | 221 CHLOROMETHANE <74-87-3> RO#2 |
| 3 | 231 VINYL CHLORIDE <75-01-4> RO#3 |
| 4 | 220 BROMOMETHANE <78-83-9> RO#4 |
| 5 | 209 CHLOROETHANE <75-00-3> RO#5 |
| 6 | 216 1,1-DICHLOROETHENE <75-35-4> RO#8 |
| 7 | 254 CARBON DISULFIDE <75-15-0> RO#9 |
| 8 | 252 ACETONE (2-PROPANONE) <67-64-1> RO#13 |
| 9 | *249 1,4-DIFLUOROBENZENE (IS) <540-36-3> RO#14 |
| 10 | 222 METHYLENE CHLORIDE <75-09-2> RO#16 |
| 11 | 226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17 |
| 12 | 214 1,1-DICHLOROETHANE <75-34-3> RO#19 |
| 13 | 257 VINYL ACETATE <108-05-4> RO#20 |
| 14 | 237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21 |
| 15 | 253 2-BUTANONE <78-93-3> RO#22 |
| 16 | 211 CHLOROFORM <67-66-2> RO#23 |
| 17 | 227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24 |
| 18 | 206 CARBON TETRACHLORIDE <56-23-5> RO#25 |
| 19 | 203 BENZENE <71-43-2> RO#26 |
| 20 | 215 1,2-DICHLOROETHANE <107-06-2> RO#27 |
| 21 | *270 D5-CHLOROBENZENE (IS) RO#29 |
| 22 | 229 TRICHLOROETHENE <79-01-6> RO#30 |
| 23 | 217 1,2-DICHLOROPROPANE <78-87-5> RO#31 |
| 24 | 212 BROMODICHLOROMETHANE <75-27-4> RO#33 |
| 25 | 218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35 |
| 26 | 256 4-METHYL-2-PENTANONE <108-01-1> RO#36 |
| 27 | 225 TOLUENE <108-88-3> RO#37 |
| 28 | 250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38 |
| 29 | 228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39 |
| 30 | 224 TETRACHLOROETHENE <127-18-4> RO#41 |
| 31 | 255 2-HEXANONE <591-78-6> RO#42 |
| 32 | 208 DIBROMOCHLOROMETHANE <124-48-1> RO#43 |
| 33 | 207 CHLOROBENZENE <108-90-7> RO#45 |
| 34 | 219 ETHYLBENZENE <100-41-4> RO#47 |
| 35 | 330 M,P-XYLENE <133-02-7> RO#48 |
| 36 | 239 O-XYLENE <133-02-7> RO#49 |
| 37 | 251 STYRENE <100-42-5> RO#50 |
| 38 | 205 BROMOFORM <75-25-2> RO#51 |
| 39 | 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54 |
| 40 | *258 D4-1,2-DICHLOROETHANE RO#57 |
| 41 | *247 BROMOFLUOROBENZENE <460-00-4> RO#58 |
| 42 | *233 D8-TOLUENE RO#59 |

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA (HGHT)	AMOUNT	%TOT
1	128	464	5.48	1	1.000	A BB	52467.	50.000 UG/KG	2.69
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	223	2:47	1	0.481	A BB	21293.	64.779 UG/KG	3.48%
9	114	610	7:37	9	1.000	A BB	198306.	50.000 UG/KG	2.69%
10	84	272	3:24	1	0.586	A BB	23058.	12.079 UG/KG	0.65%
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	446	5:34	1	0.961	A BB	6558.	46.449 UG/KG	2.50%
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	544	6:48	9	0.892	A BB	99903.	25.277 UG/KG	1.36%
20	62	NOT FOUND							
21	117	1001	12:31	21	1.000	A BB	195114.	50.000 UG/KG	2.69%
22	130	NOT FOUND							
23	63	658	8:13	9	1.079	A BB	1872.	1.333 UG/KG	0.07%
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	801	10:01	21	0.800	A BB	17674.	12.528 UG/KG	0.67%
27	92	809	10:07	21	0.808	A BB	915382.	279.559 UG/KG	15.03%
28	75	NOT FOUND							
29	97	NOT FOUND							
30	164	NOT FOUND							
31	43	920	11:30	21	0.919	A VB	306556.	272.924 UG/KG	14.67%
32	129	NOT FOUND							
33	112	NOT FOUND							
34	106	1030	12:52	21	1.029	A BV	369299.	171.818 UG/KG	9.24%
35	106	1049	13:07	21	1.048	A VB	1079840.	337.204 UG/KG	18.13%
36	106	1104	13:48	21	1.103	A BB	714692.	249.421 UG/KG	13.41%
37	104	1108	13:51	21	1.107	A BB	492304.	95.283 UG/KG	5.12%
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	540	6:45	1	1.164	A BB	81903.	47.219 UG/KG	2.54%
41	95	1182	14:46	21	1.181	A BB	164551.	49.670 UG/KG	2.67%
42	98	800	10:00	21	0.799	A BB	203086.	44.872 UG/KG	2.41%

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:55	0.98	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:07		10.000			50.00		0.930	
3	1:13		10.000			50.00		1.260	
4	1:31		10.000			50.00		1.632	
5	1:37		10.000			50.00		0.710	
6	2:40		5.000			50.00		1.612	
7	2:50		5.000			50.00		4.085	
8	2:52	0.97	10.000	0.05	64.78	50.00	0.406	0.313	1.30
9	7:42	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	3:28	0.98	5.000	0.12	12.08	50.00	0.439	1.819	0.24
11	3:54		5.000			50.00		1.643	
12	4:36		5.000			50.00		2.402	
13	4:52		10.000			50.00		0.538	
14	5:34		5.000			50.00		1.673	

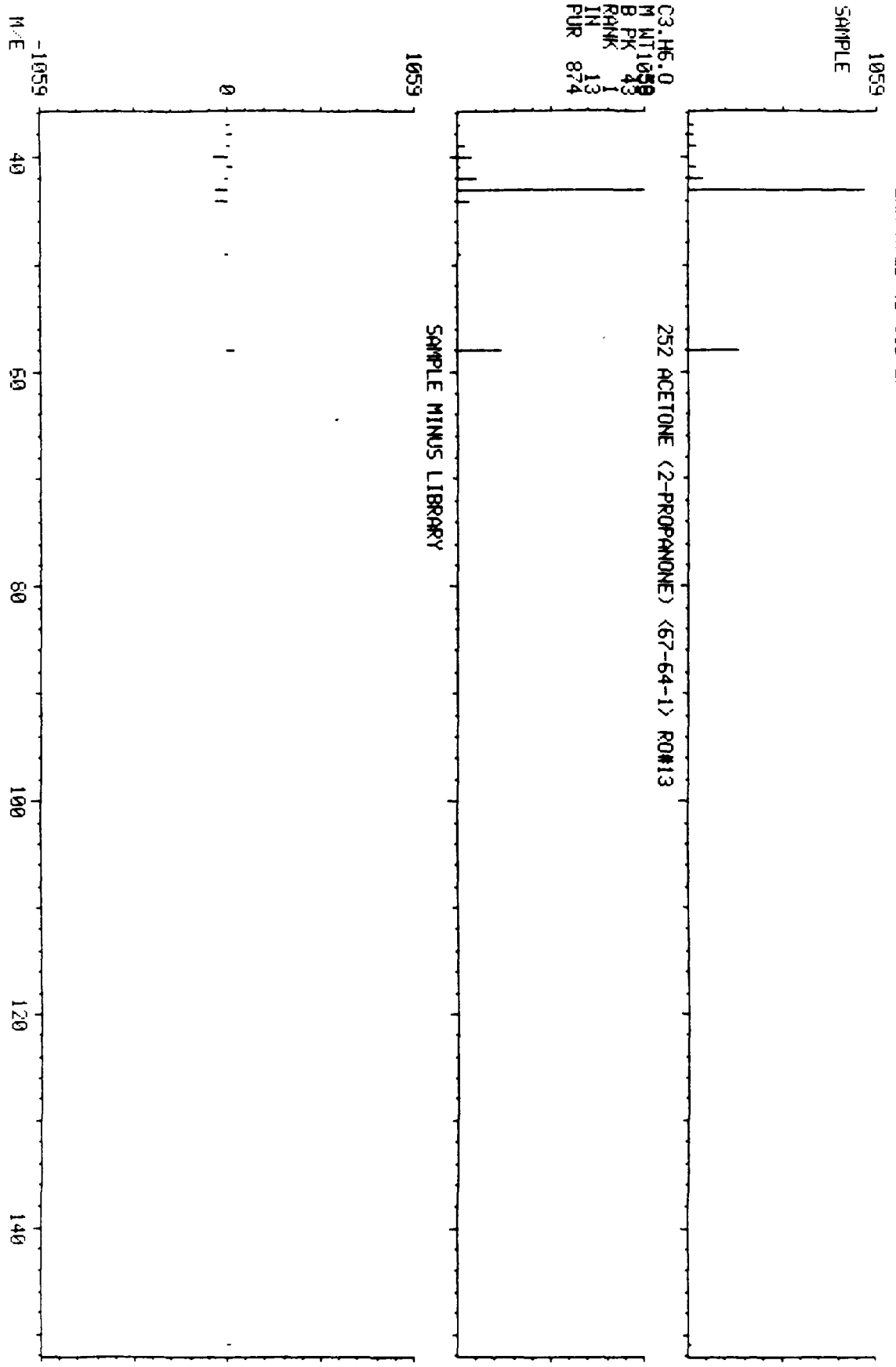
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:41	0.98	10.000	0.10	46.45	50.00	0.125	0.135	0.93
16	6:10		5.000			50.00		3.249	
17	6:19		5.000			50.00		0.880	
18	6:34		5.000			50.00		0.898	
19	6:53	0.99	5.000	0.18	25.28	50.00	0.504	0.997	0.51
20	6:58		5.000			50.00		2.119	
21	12:34	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	8:01		5.000			50.00		0.581	
23	8:21	0.99	5.000	0.22	1.33	50.00	0.009	0.354	0.03
24	8:55		5.000			50.00		0.785	
25	9:41		5.000			50.00		0.771	
26	10:04	0.99	15.000	0.05	12.53	50.00	0.091	0.362	0.25
27	10:10	0.99	5.000	0.16	279.56	50.00	4.692	0.839	5.59
28	10:41		5.000			50.00		0.357	
29	10:57		5.000			50.00		0.407	
30	11:06		5.000			50.00		0.673	
31	11:34	0.99	15.000	0.06	272.92	50.00	1.571	0.288	5.46
32	11:34		5.000			50.00		0.776	
33	12:37		5.000			50.00		1.212	
34	12:55	1.00	5.000	0.21	171.82	50.00	1.893	0.551	3.44
35	13:09	1.00	5.000	0.21	337.20	50.00	5.534	0.821	6.74
36	13:51	1.00	5.000	0.22	249.42	50.00	3.663	0.734	4.99
37	13:54	1.00	5.000	0.22	95.28	50.00	2.523	1.324	1.91
38	14:07		5.000			50.00		0.577	
39	15:15		5.000			50.00		0.655	
40	6:51	0.99	5.000	0.23	47.22	50.00	1.561	1.653	0.94
41	14:49	1.00	5.000	0.24	49.67	50.00	0.843	0.849	0.99
42	10:04	0.99	5.000	0.16	44.87	50.00	1.041	1.160	0.90

COMPUchem LABS
LIBRARY SEARCH
12/27/89 5:39:00 + 2:47
SAMPLE: 16 IC#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (5 158 2N 0T)

DATA: GR009686C19 # 223

BASE M/E: 43
RID: 4105.

C3.H6.0
M.WT 1059
B.PK 43
RANK 1
IN 13
PUR 874

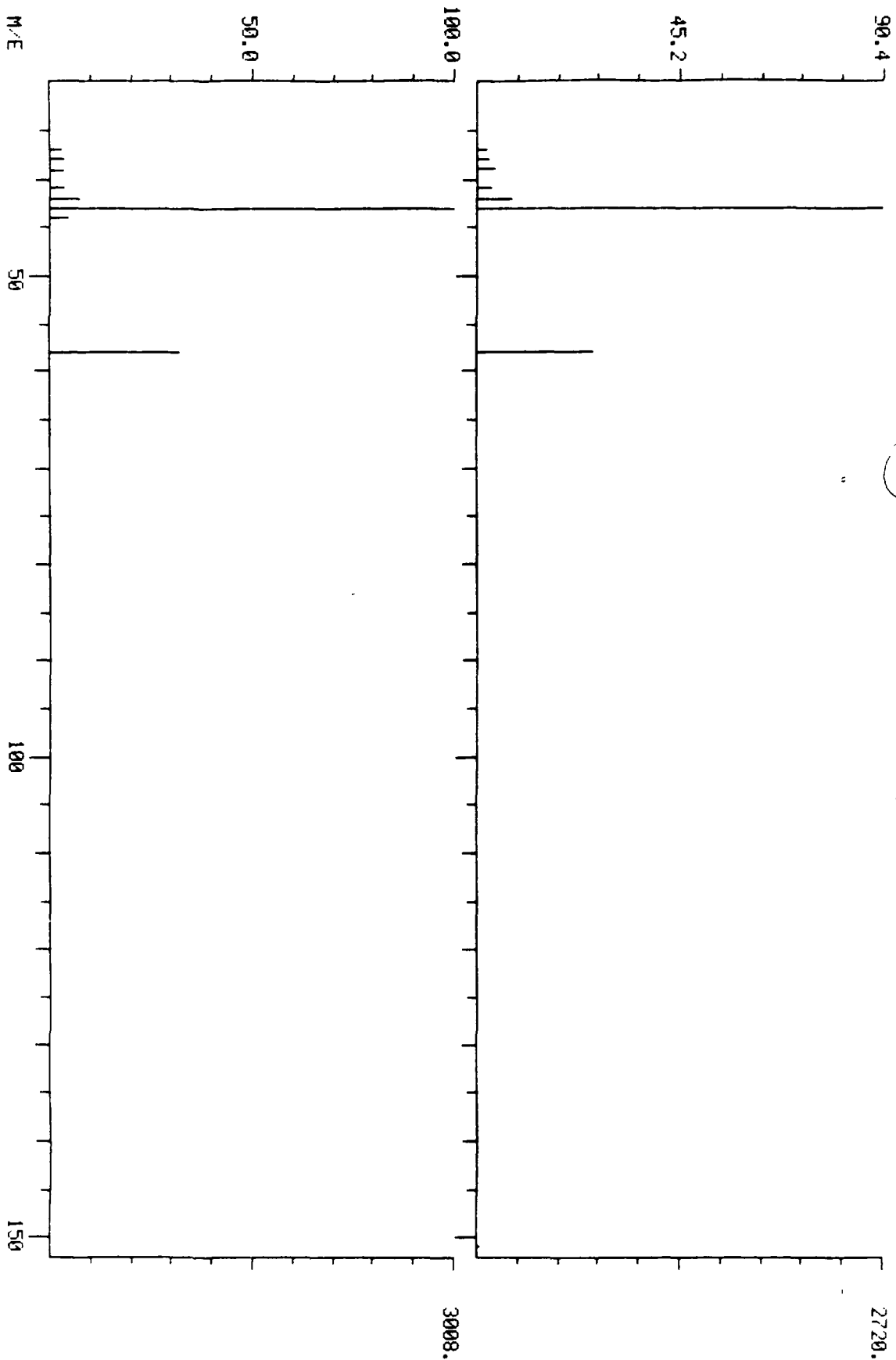


DUAL MASS SPECTRUM
12/27/89 5:39:00 + 2:47
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (5 15B 2M) (252) ACETONE (2-PROPANONE) <67-64-1> R0#13

COMPUchem LABS

DATA: GR009686C19 #228

BASE M/E: 43/ 43
RIC: 4103. / 4695.



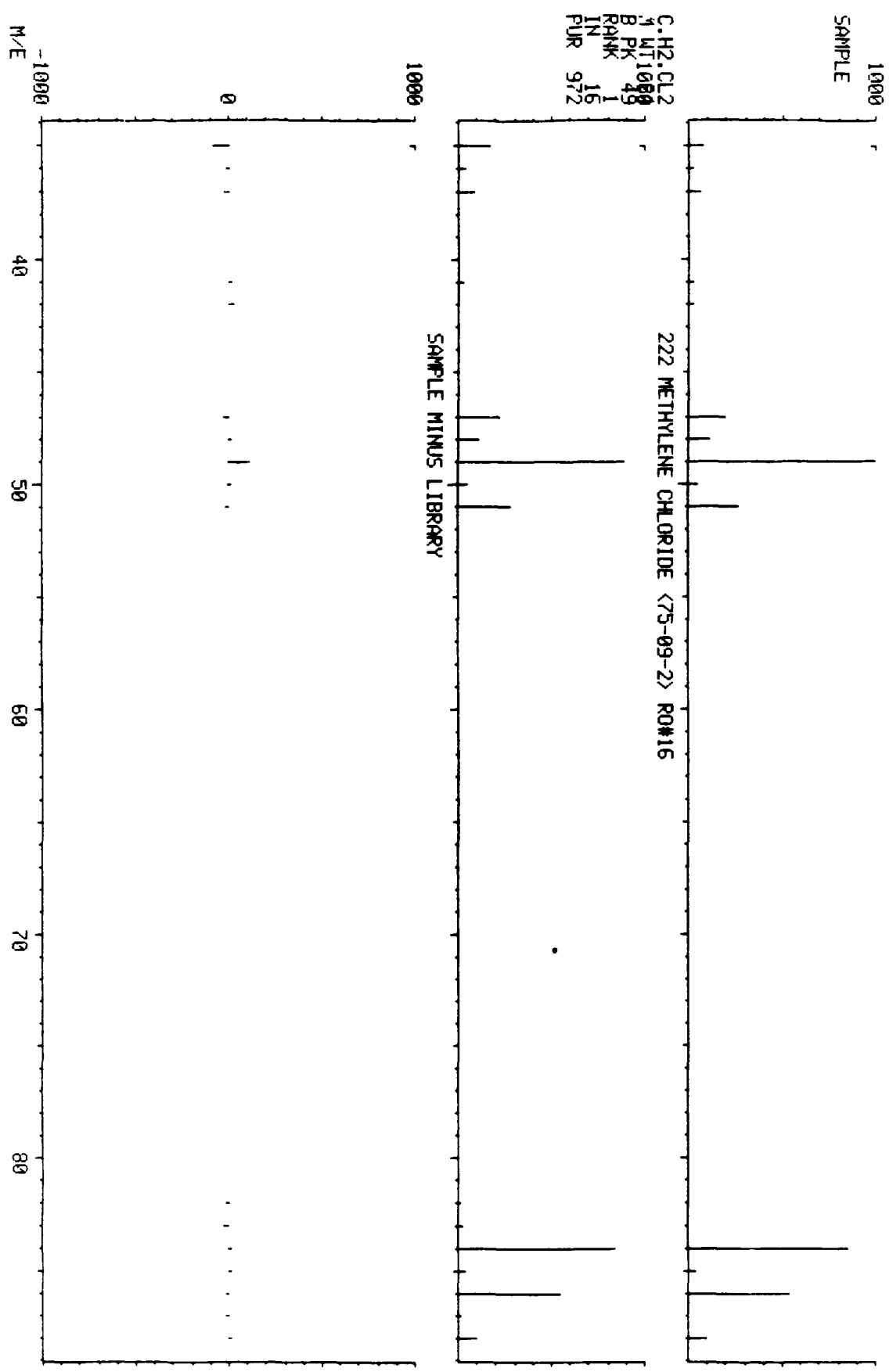
LIBRARY SEARCH
12/27/89 5:39:00 + 3:24
SAMPLE: 16 CC#S09686 CASE#18756.7 EPA#B2018 QN#19
ENHANCED (S 158 2N 0T)

COMPUCHEM L865

DATA: GR005686C19 # 272

BASE M/E: 49
RIC: 11711.

C-H2.C12
Y UT 1000
B PK 49
RANK 1
IN 16
PUR 972

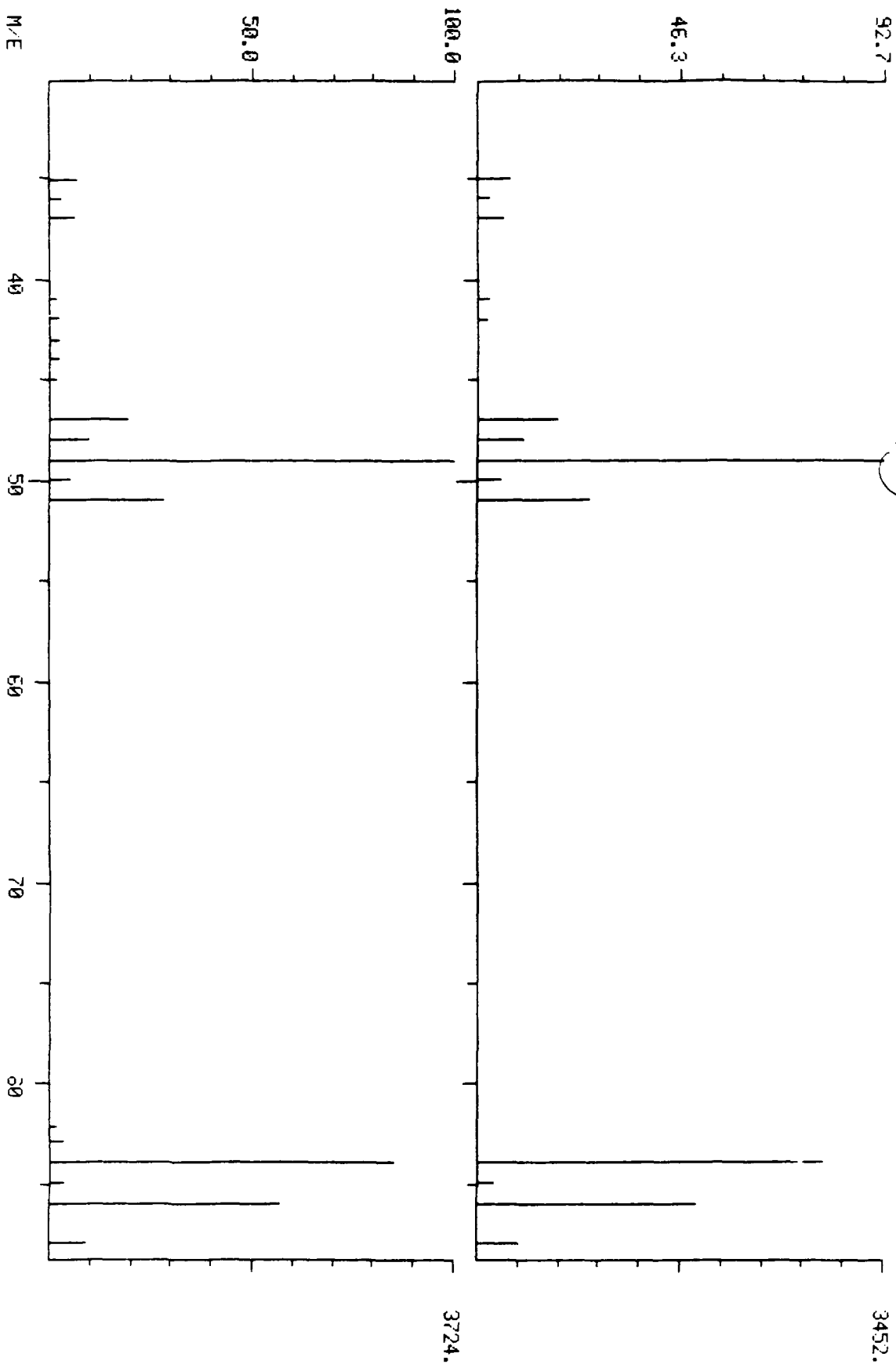


DUAL MASS SPECTRUM
12/27/89 5:39:00 + 3:24
SAMPLE: 1G CC#309586 CASE#18756.7 EPA#B2018 ON#19
ENHANCED (S 15B 2N) 222 METHYLENE CHLORIDE (75-09-2) RQ#16

COMFUCHEM LABS

DATA: GR009586C19 #272

BASE M/E: 49/ 49
RIC: 11711.7 12863.



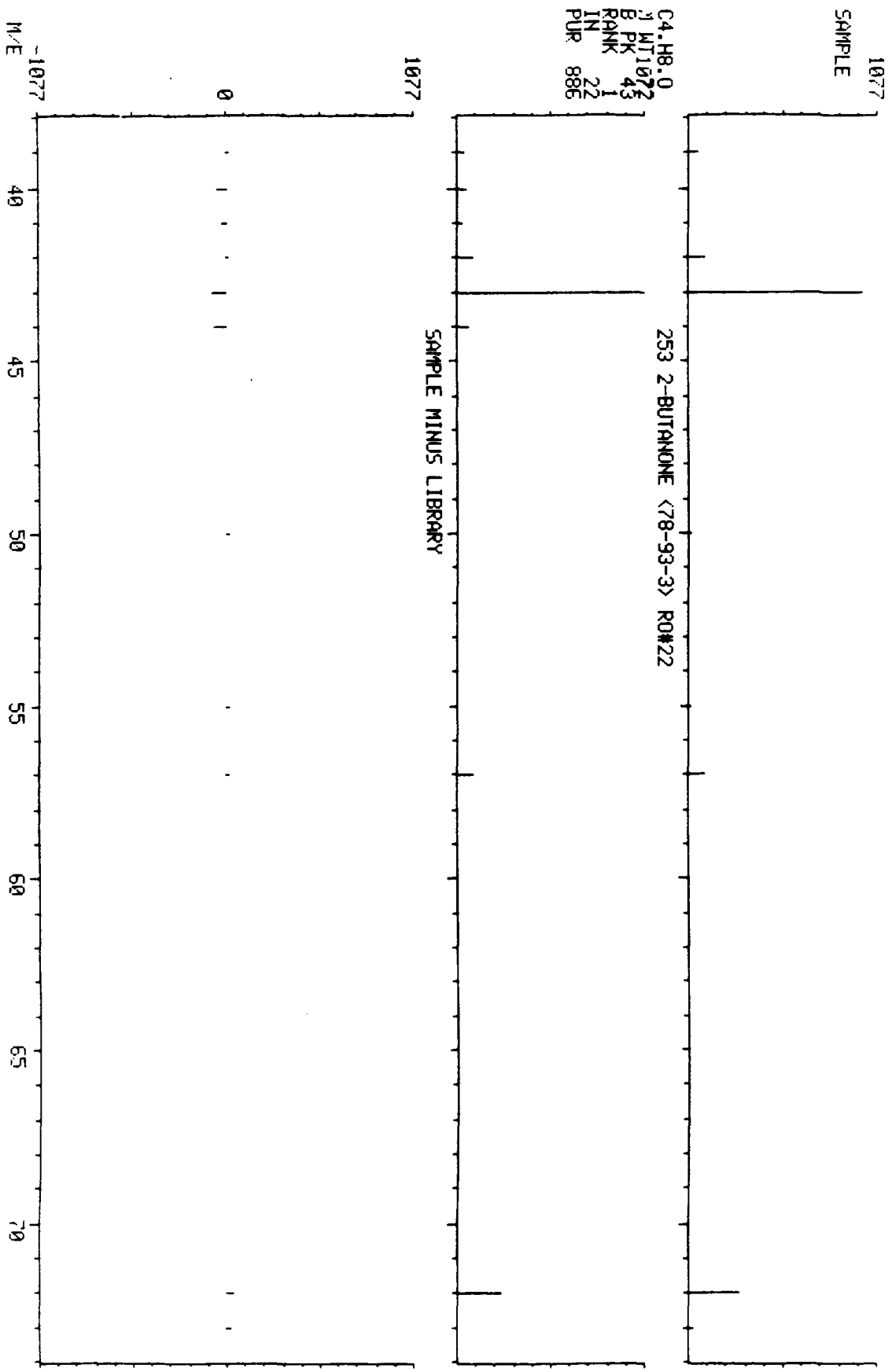
LIBRARY SEARCH
12/27/89 5:39:00 + 5:34
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B2018 ON#19
ENHANCED (5 158 2N 01)

COMPUCHEN LABS

DATA: GR009686C19 # 446

BASE M/E: 43
RIC: 2959.

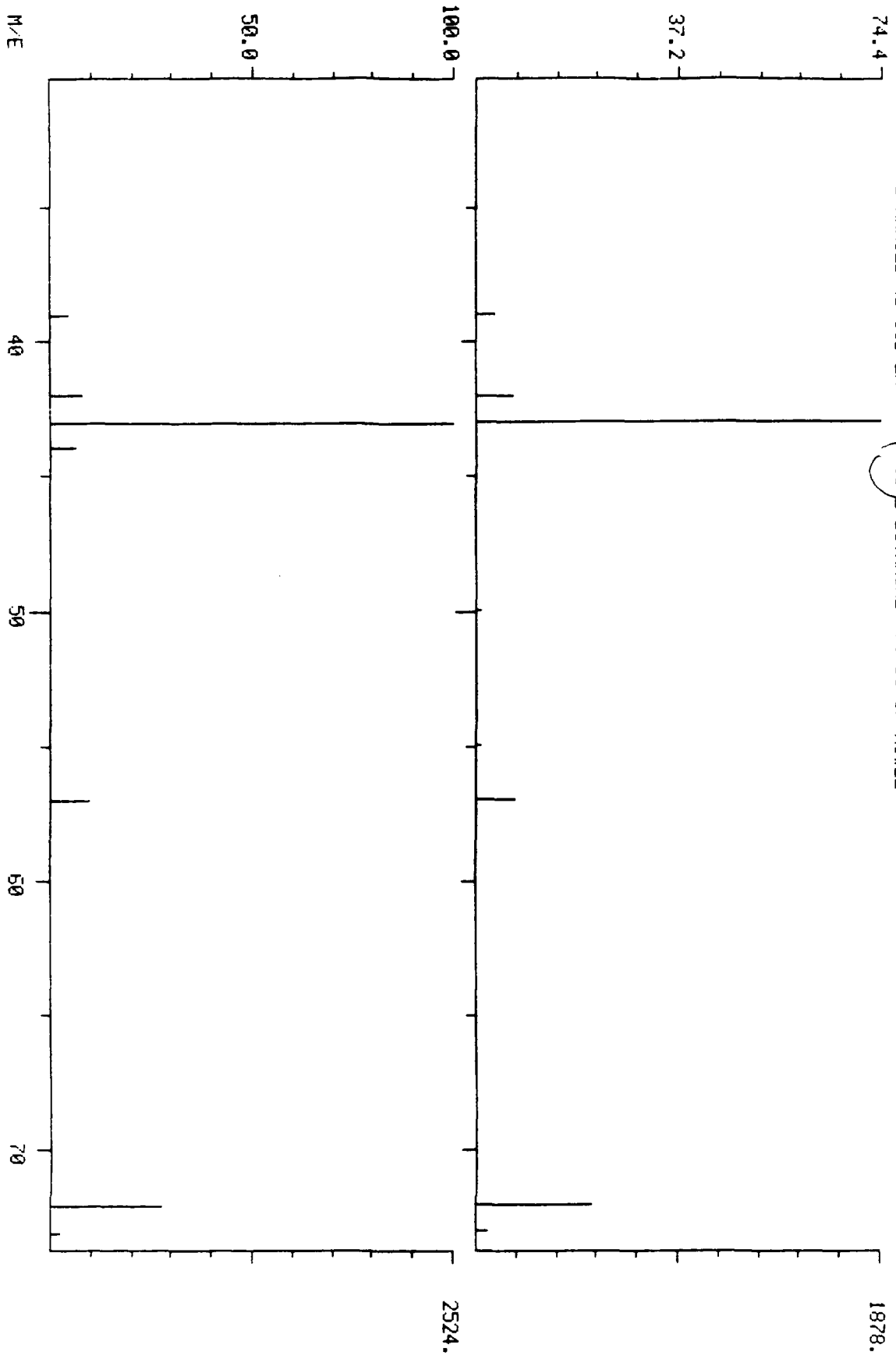
C4.H8.0
M WT 1077
B PK 43
RANK 1
IN 22
PUR 886



DUAL MASS SPECTRUM
12/27/89 5:39:00 + 5:34
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (5 158 2N) 253 2-BUTANONE (78-93-3) RM#22

COMFUCHEM LABS

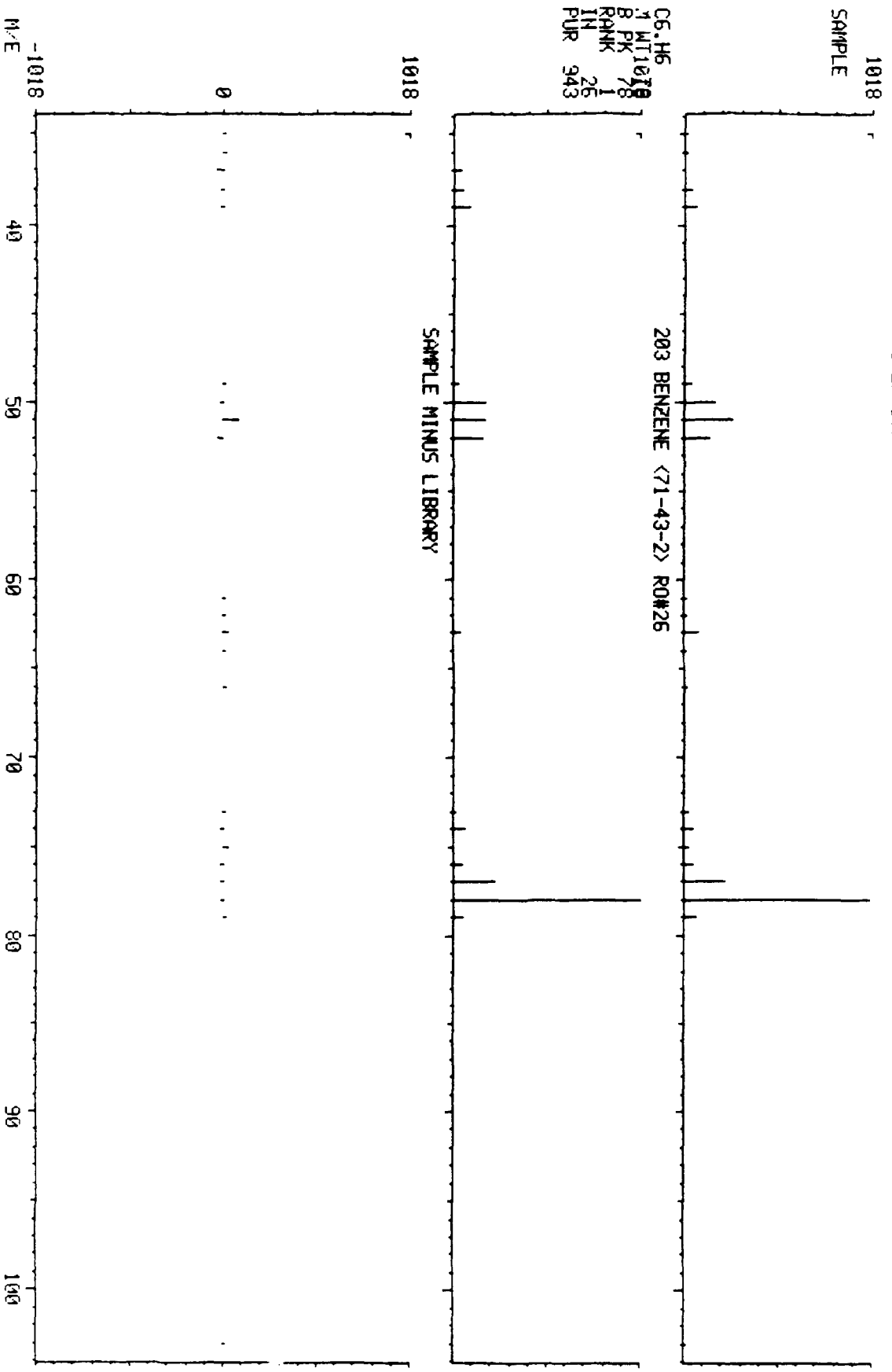
DATA: GR009686019 #446 BASE M/E: 43/ 45
RIC: 2959.7 3999.



COMPUCHEN LABS
LIBRARY SEARCH
12/27/89 5:39:00 + 6:42
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B DN#19
ENHANCED (5 15B 2N 0T)

DATA: GR009686C19 # 544
BASE M/E: 78
RIC: 21503.

CG.H6
M WT 1070
B PK 78
RANK 1
IN 26
PUR 943

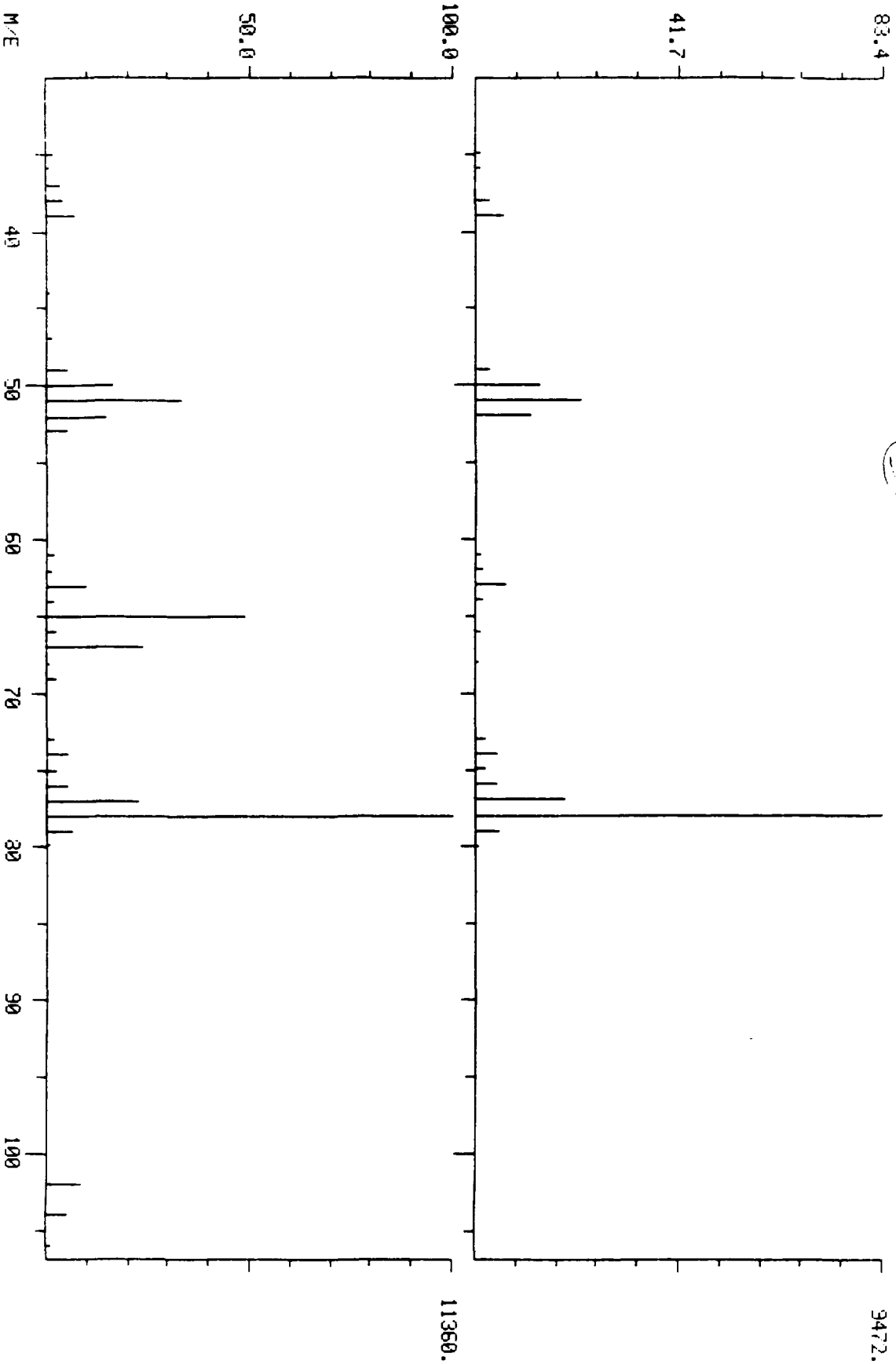


CUHL MASS SPECTRUM
12/27/89 5:39:00 + 6:48
SAMPLE: 1G.CC#309686 CASE#18756.7 EPA#B2018 OH#19
ENHANCED (5 158.2H) 203/BENZENE (71-43-2) RM#26

COMPUCHEN LABS

DATA: GR009686C19 #544

BASE M/E: 78/ 78
RIC: 21599. / 58911.



COMPUchem LABS
LIBRARY SEARCH
12/27/89 5:39:00 + 10:01
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (S 158 ZN 0T)

DATA: GR009686C19 # 801

BASE M/E: 98
RIC: 45783.

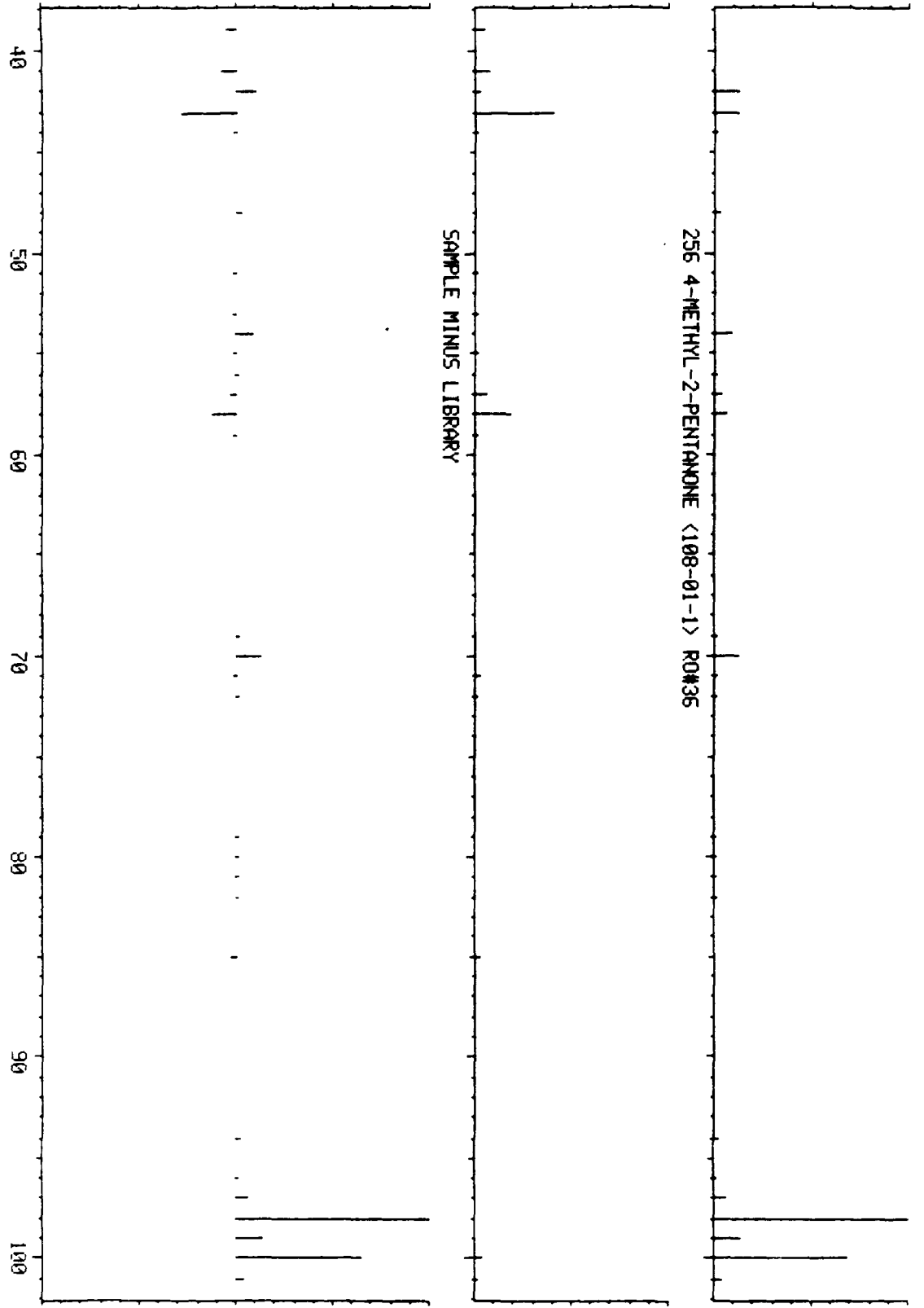
1000
SAMPLE

CG.H12.0
Y.MT1000
B.PK 43
RANK 1
IN 35
PUR 136

256 4-METHYL-2-PENTANONE <108-01-1> RO#35

SAMPLE MINUS LIBRARY

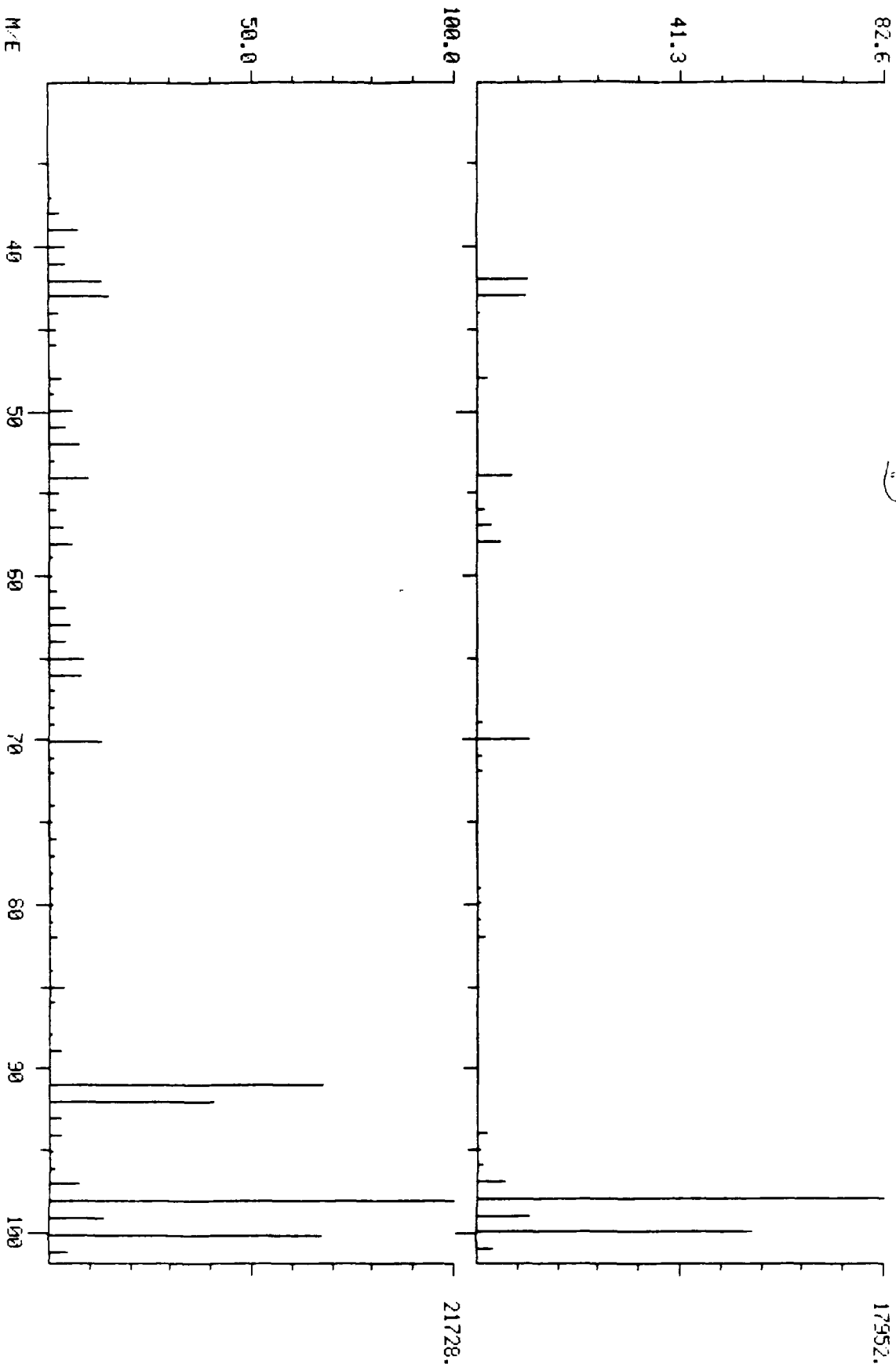
-1000
M/E



DUAL MASS SPECTRUM
12/27/89 5:39:00 + 10:01
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#82018 ON#19
ENHANCED (5 158 2N) (256) 4-METHYL-2-PENTANONE (108-01-1) R0#36

COMPUCHEN LABS

DATA: GR009686C19 #801 BASE M/E: 98/ 98
RIC: 46763.7 102527.

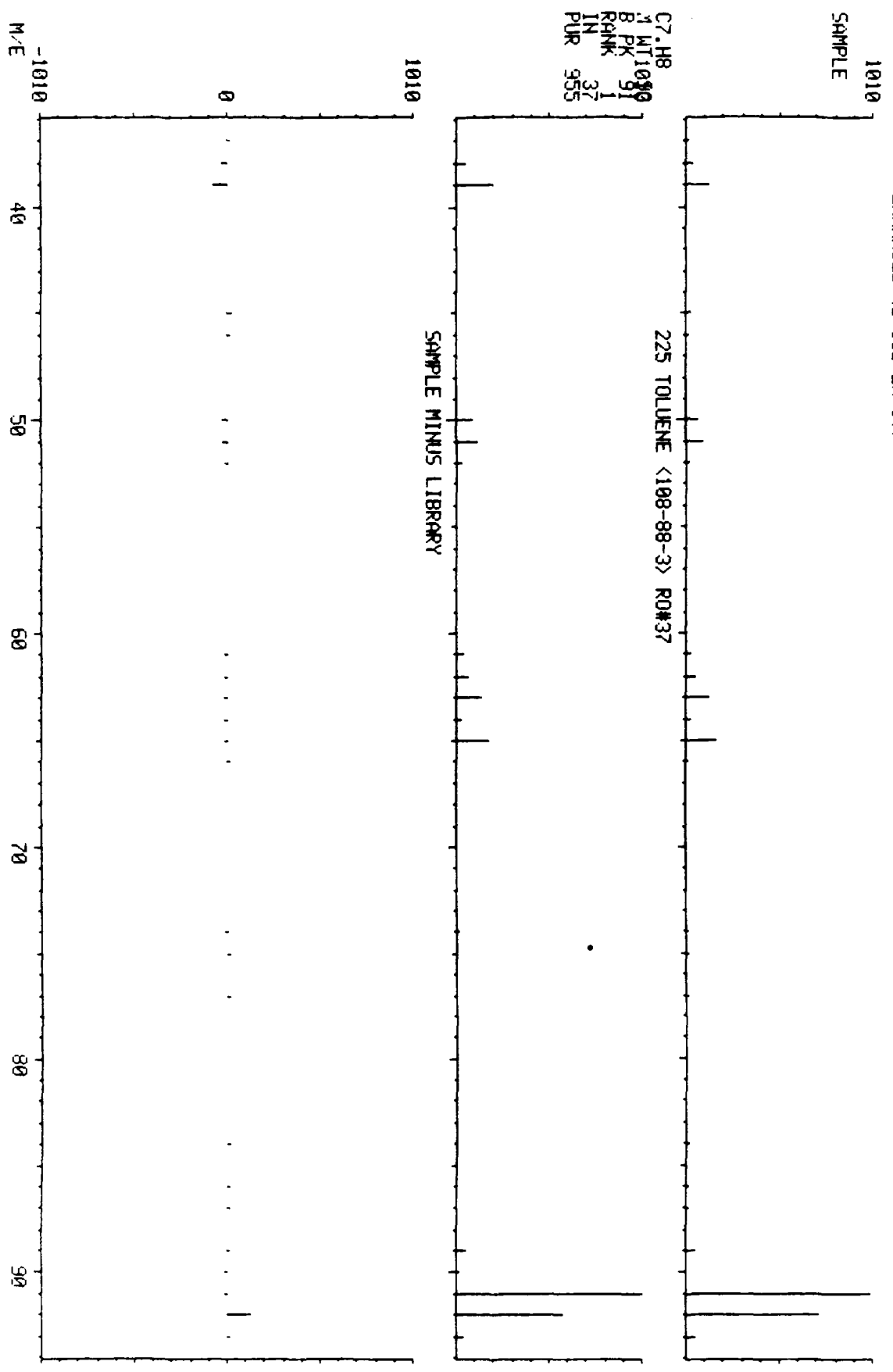


COMPUCHEM LABS
LIBRARY SEARCH
12/27/89 5:39:00 + 10:07
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B QM#19
ENHANCED (5 15B 2N 0T)

DATA: GP009686C19 # 809

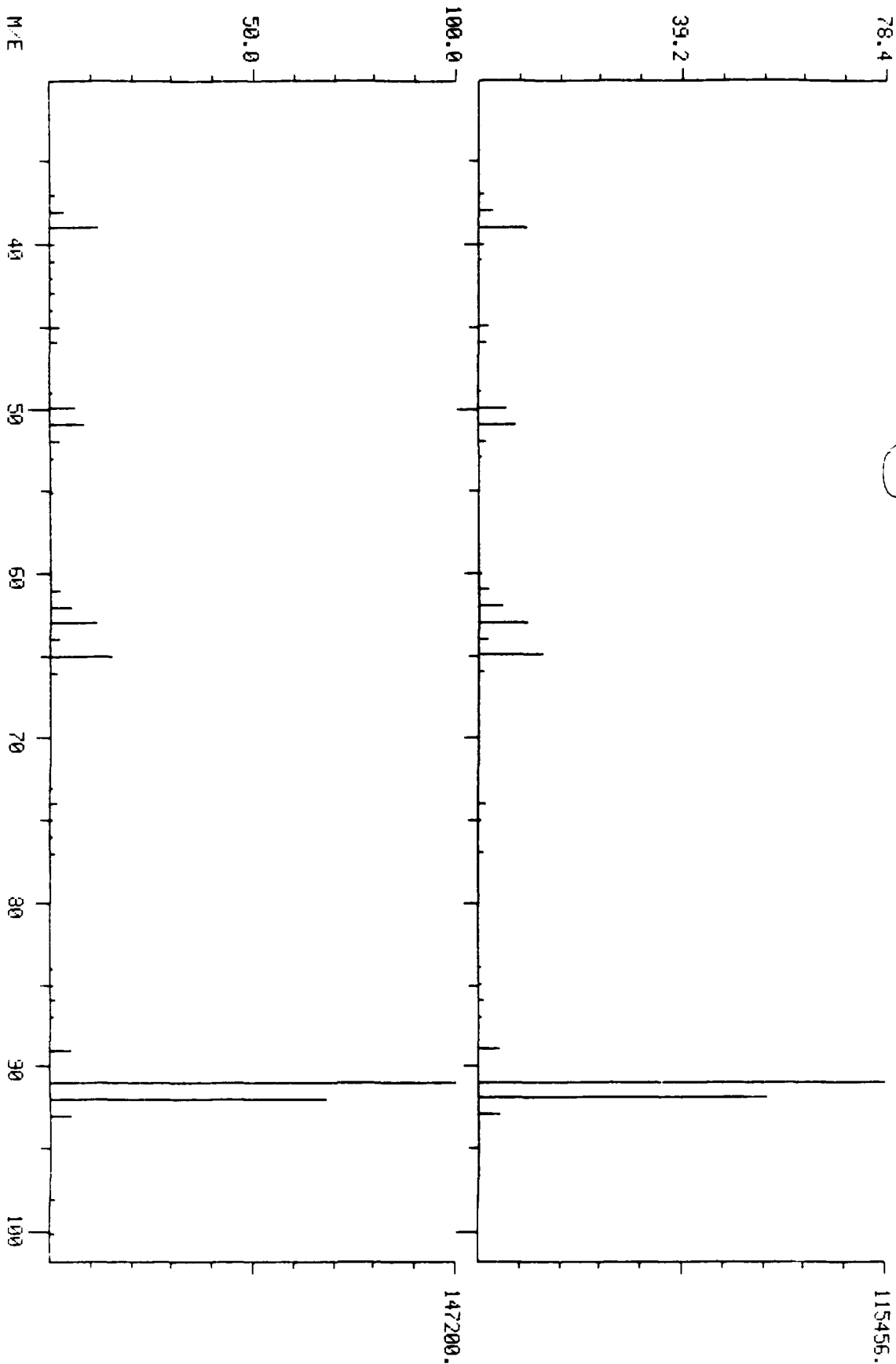
BASE M/E: 91
RIC: 204639.

C7.H8
1 MT 1090
8 PK 91
RANK IN 37
PUR 955



DUAL MASS SPECTRUM
12/27/89 5:39:00 + 10:07
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#R2018 QN#19
ENHANCED (5 158 2N) 225 TOLUENE <106-88-3> RM#37

COMPUchem LABS
DATA: GR009686C19 #809 BASE M/E: 91/ 91
RIC: 309247./ 358335.

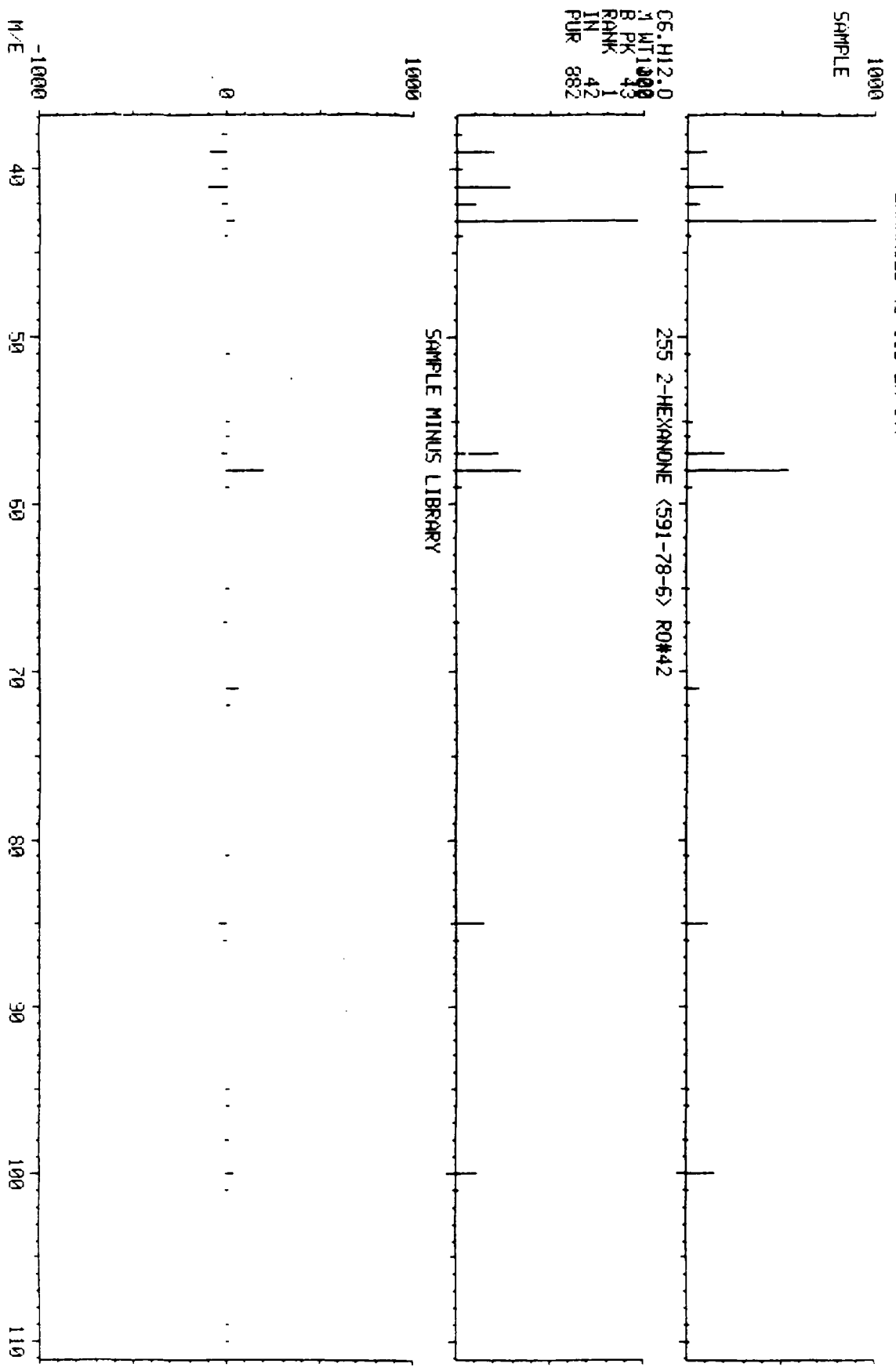


COMPUchem LABS
LIBRARY SEARCH
12/27/89 5:39:00 + 11:30
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#R2018 ON#19
ENHANCED (5.158 2H 0T)

DATA: GR009686C19 # 920

BASE M/E: 43
PIC: 77439.

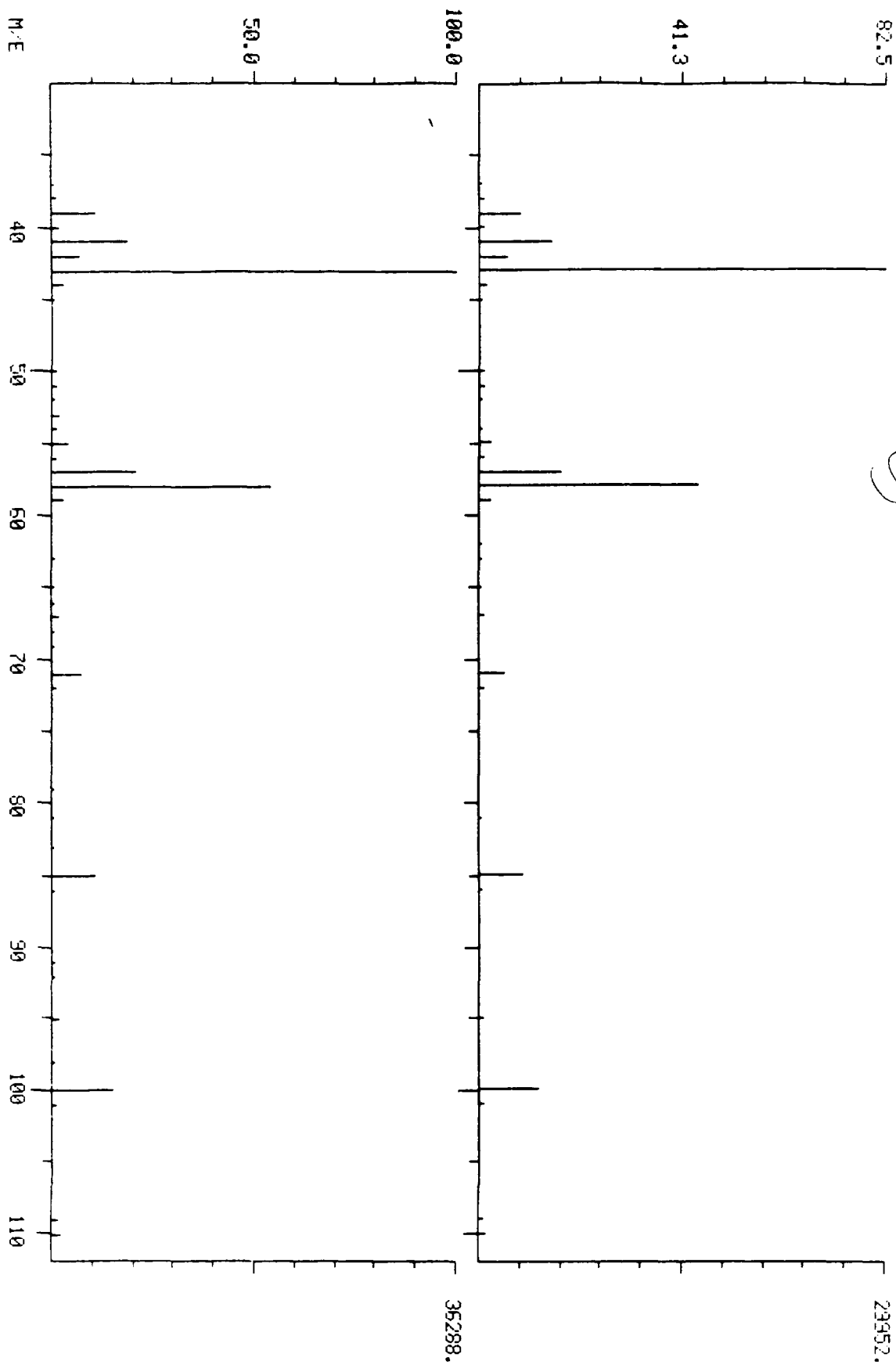
06.H12.0
I: WT 1000
R PK 43
RANK 42
IN 1
PUR 882



DUAL MASS SPECTRUM
12/27/89 5:39:00 + 11:30
SAMPLE: 16 CC#309686 CASE#18756.7 EPA#82018 QN#19
ENHANCED (5 158 2N) (255) 2-HEXANONE (591-78-6) R0#42

COMPUCHEM LABS

DATA: GR009686C19 #920 BASE M/E: 43/ 43
RIC: 79859.7 102271.

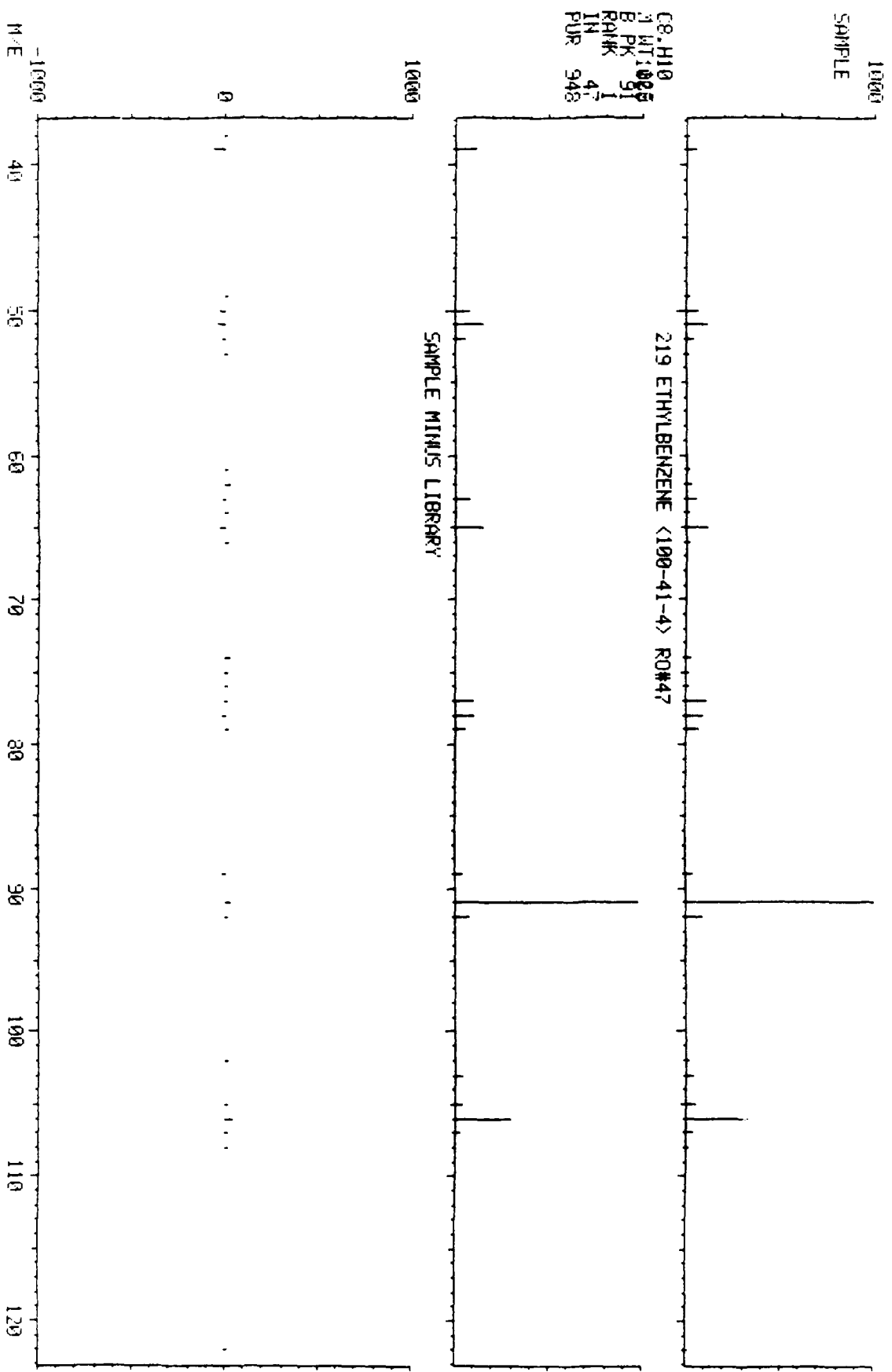


COMPUchem LABS
LIBRARY SEARCH
12/27/89 5:39:00 + 12:52
SAMPLE: 1G.CC#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (5 158 2H 0T)

DATA: GR009686C19 #1030

BASE M/E: 91
RIC: 247551.

CE: H10
1 M/I: 025
B PK: 91
RANK: 1
IN: 47
PUR: 948

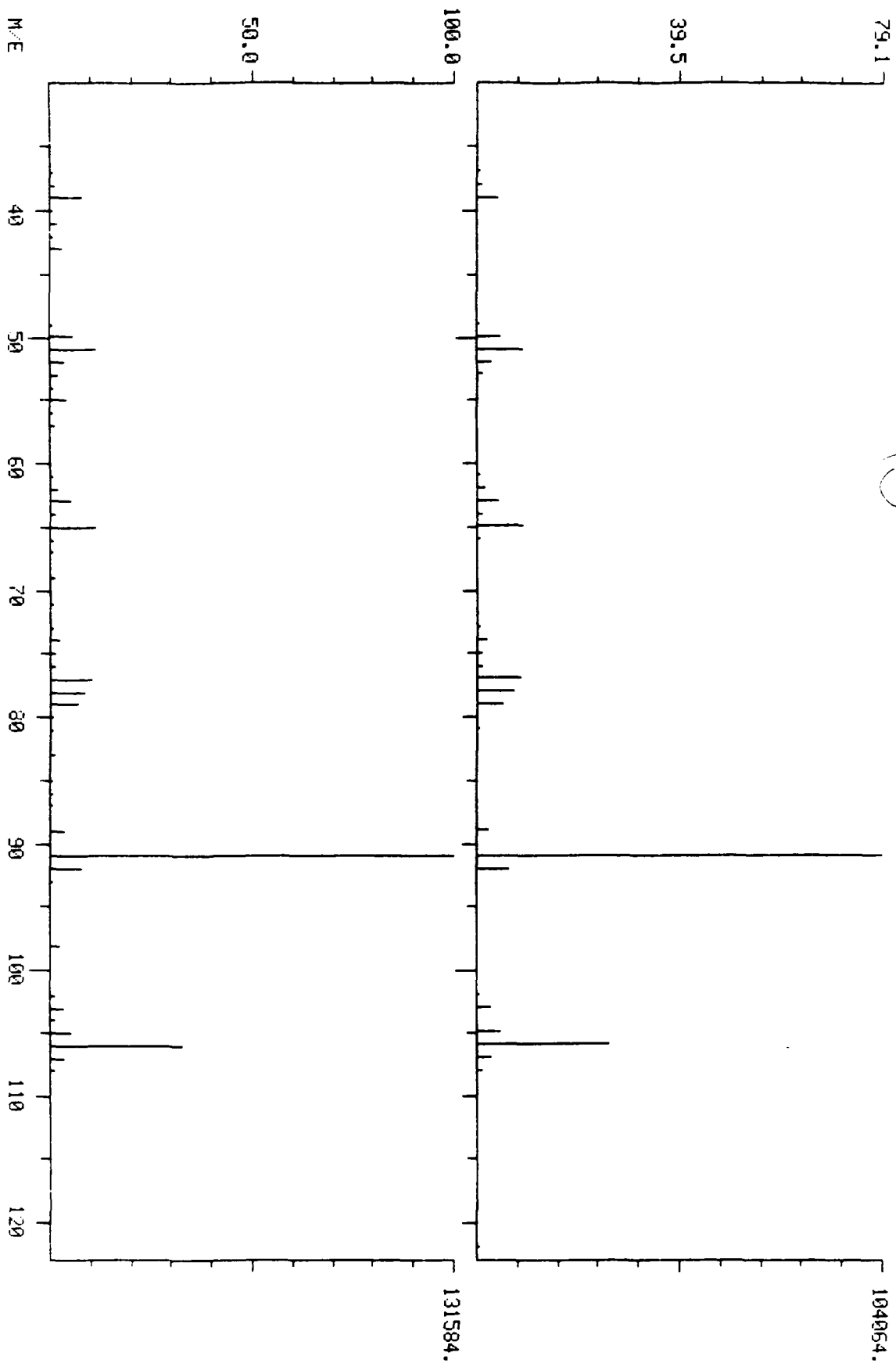


COMPUCHEM LABS

DATA: GR009686C19 #1030 BASE M/E: 91 / 91

RIC: 249599. / 348671.

DUAL MASS SPECTRUM
12/27/89 5:39:00 + 12:52
SAMPLE: 1G OC#309686 CASE#18756.7 EPA#B2018 ON#19
ENHANCED (5 15R 2H) (219 ETHYLBENZENE <100-41-4> R0#47

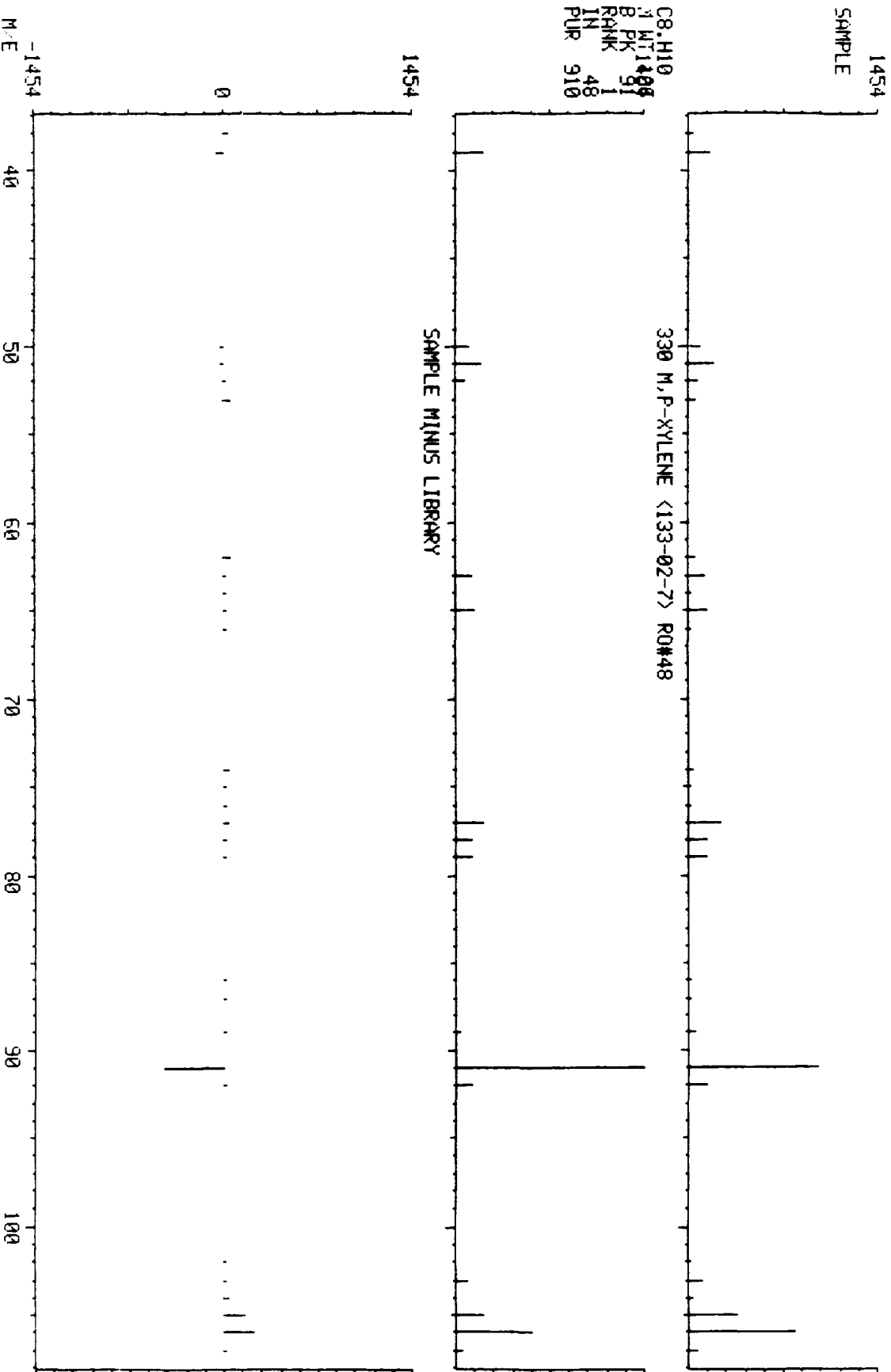


LIBRARY SEARCH
12/27/89 5:39:00 + 13:07
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B DN#19
ENHANCED (S 158 ZN 0T)

COMPUCHEM LABS

DATA: GR009686C19 #1049

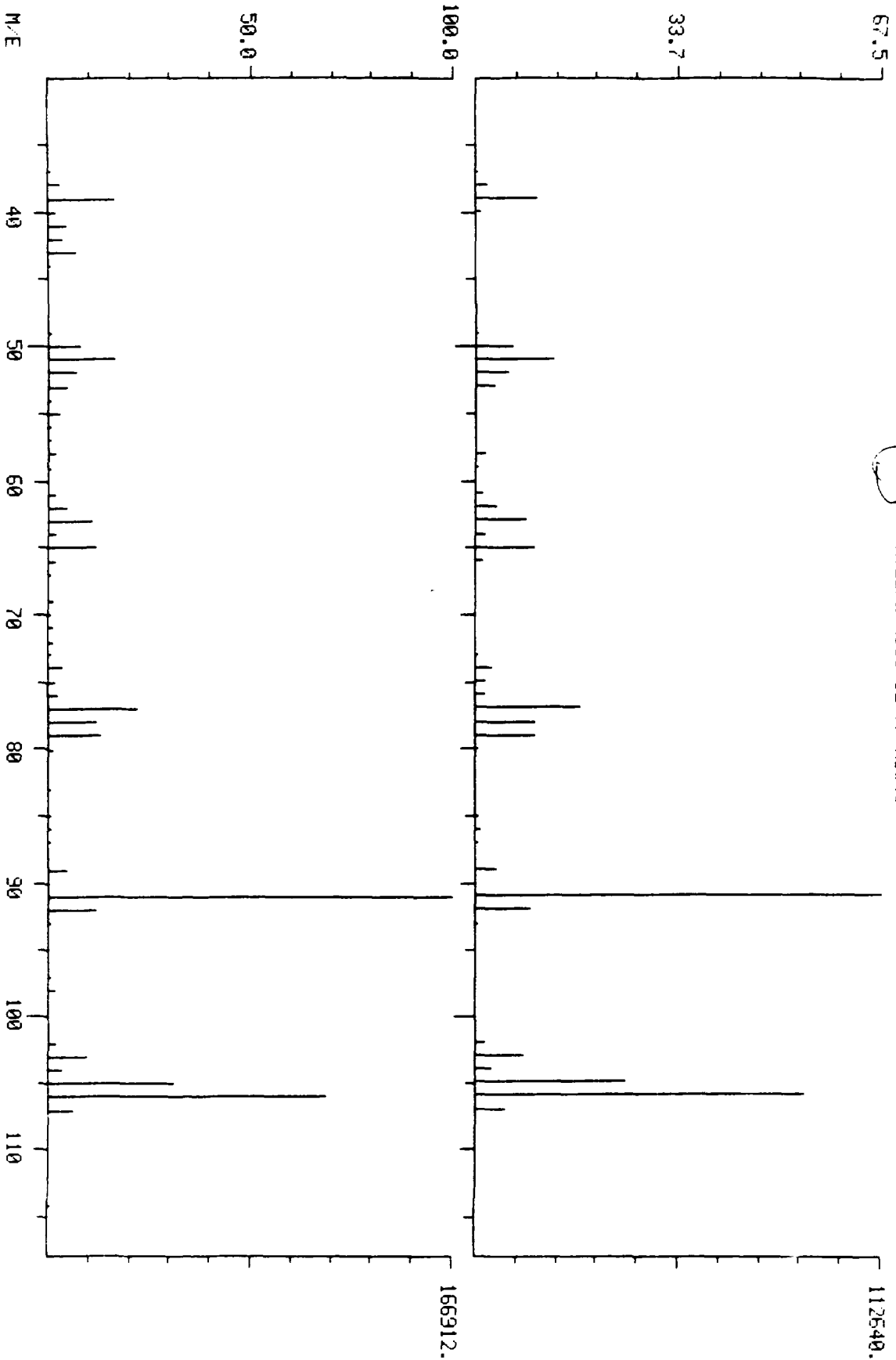
BASE M/E: 91
RIC: 475647.



DUAL MASS SPECTRUM
12/27/89 5:39:00 + 15:07
SAMPLE: 1G CCM309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (5.158 ZIN) 330 M/P-XYLENE (133-02-7) R0#48

COMPUchem LABS

DATA: GR009686C19 #1049 BASE M/E: 91 / 31
RIC: 486399. / 696319.



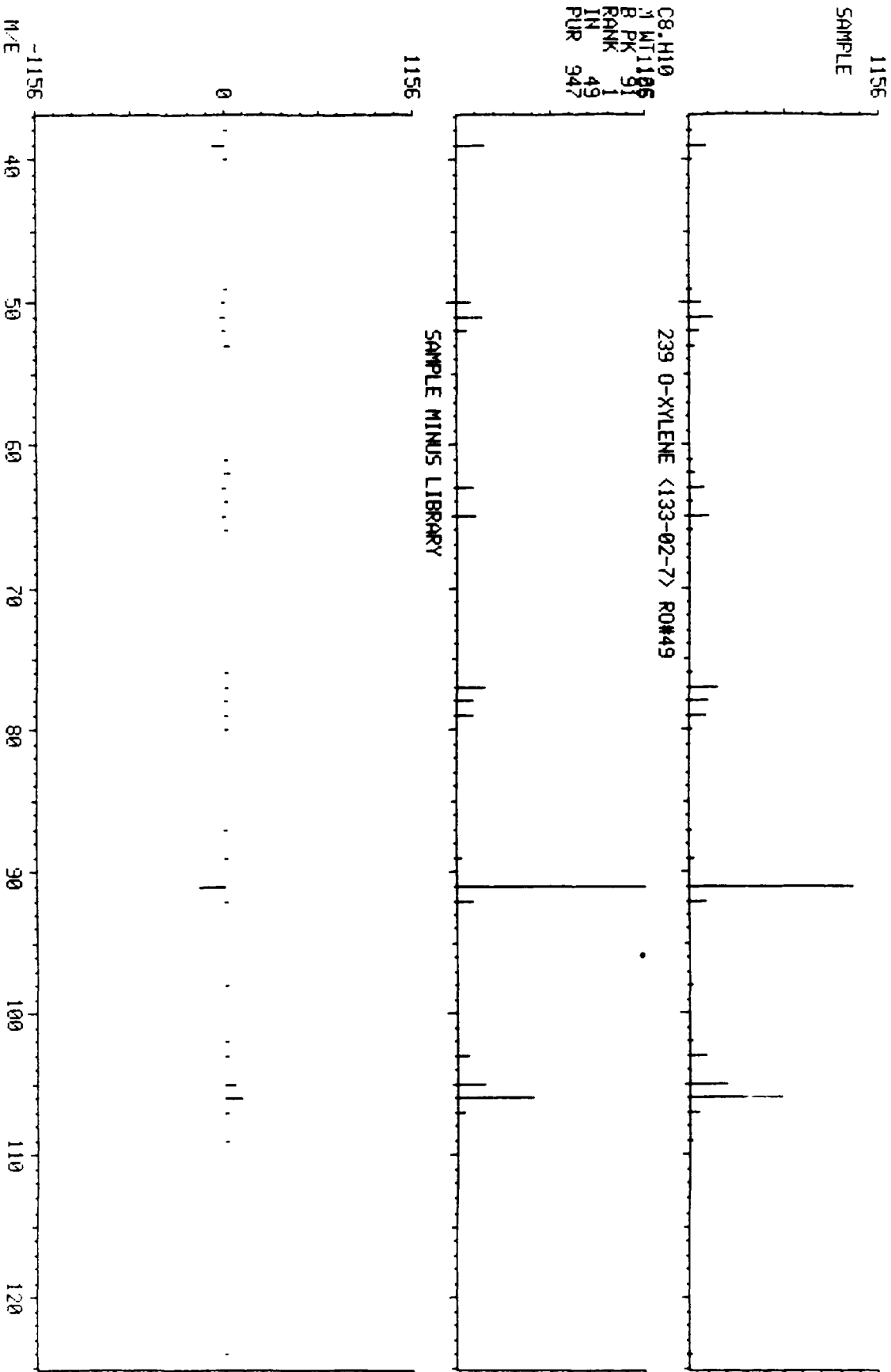
COMPUchem LABS
LIBRARY SEARCH
12/27/89 5:39:00 + 13:48
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (5 158 2N 0T)

COMPUchem LABS

DATA: GR009686C19 #1104

BASE M/E: 91
RIC: 342527.

C8.H10
Y MT 1185
B PK 91
RANK 1
IN 49
PUR 947

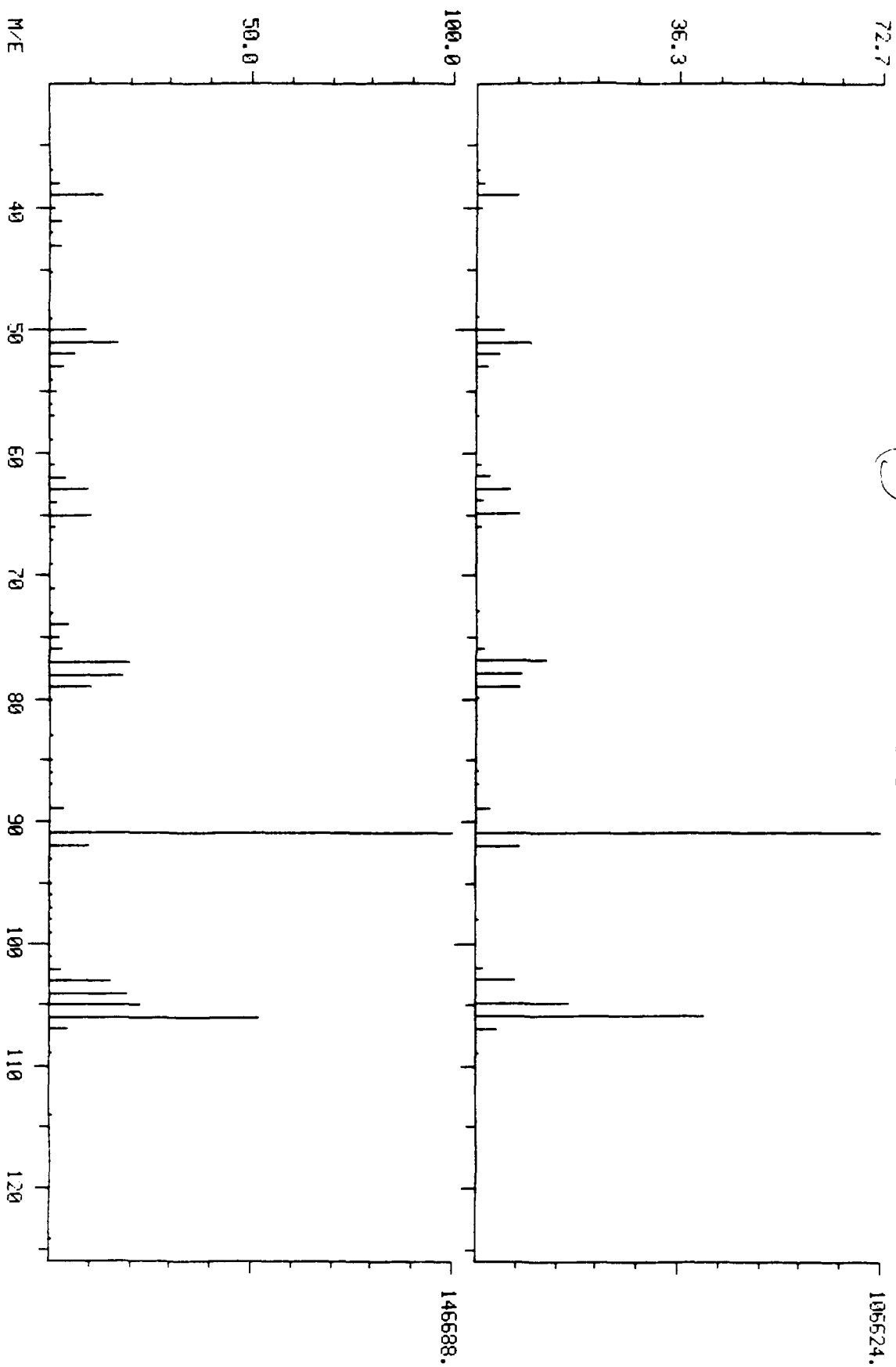


DUAL MASS SPECTRUM
12/27/89 5:39:00 + 13:48
SAMPLE: 1G CCA#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (5 15B 2N) 239 0-XYLENE <133-02-7> R0#49

COMPUCHEN LABS

DATA: GR009686C19 #1104 BASE M/E: 91 / 91

RIC: 345599. / 570367.



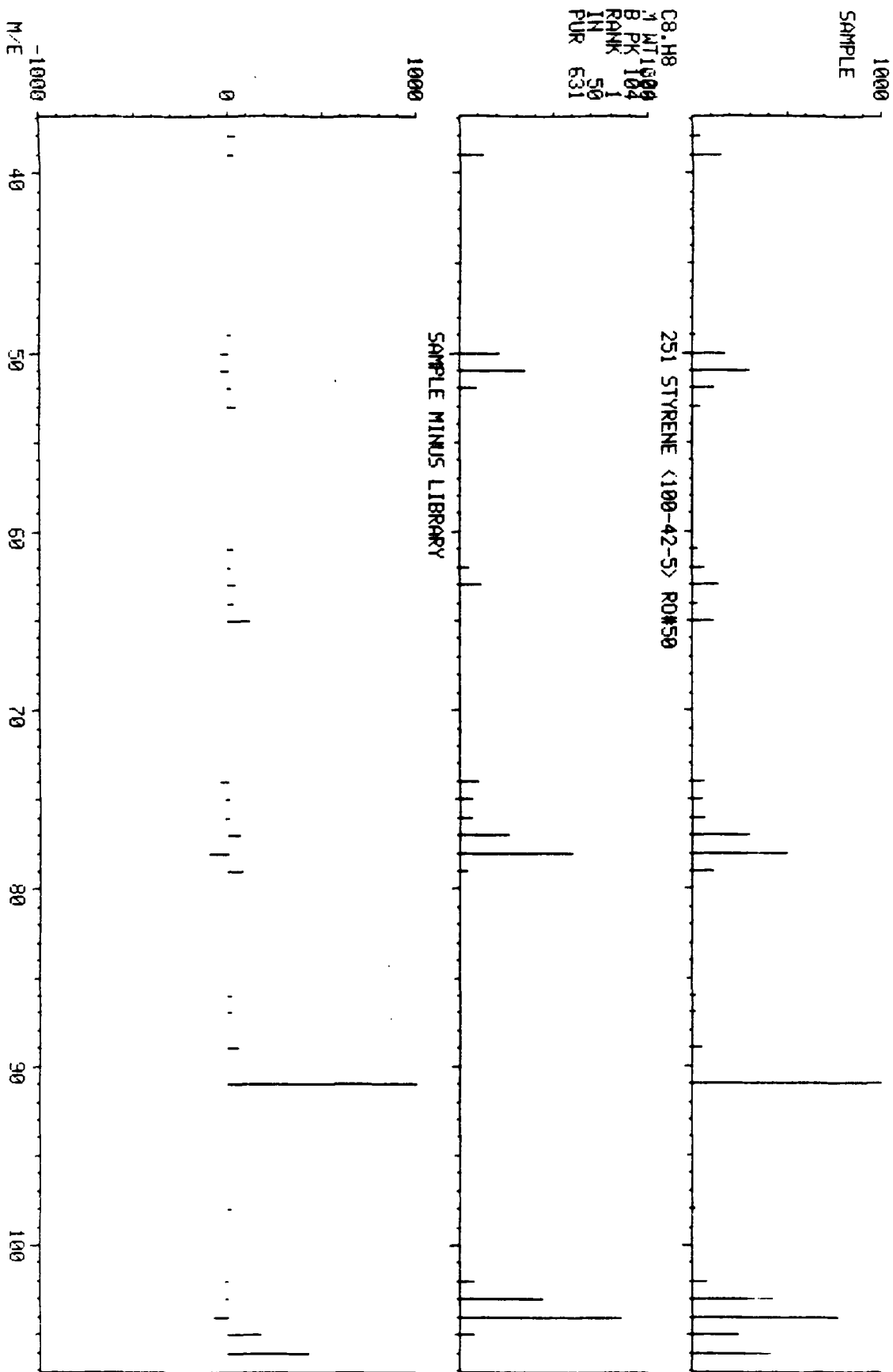
LIBRARY SEARCH
12/27/89 5:39:00 + 13:51
SAMPLE: 1G.CC#309686.CASE#18756.7 EPA#B201B QN#19
ENHANCED (S 158 2H 0T)

COMPUchem LABS

DATA: GR009686C19 #1108

BASE M/E: 91
RIC: 314879.

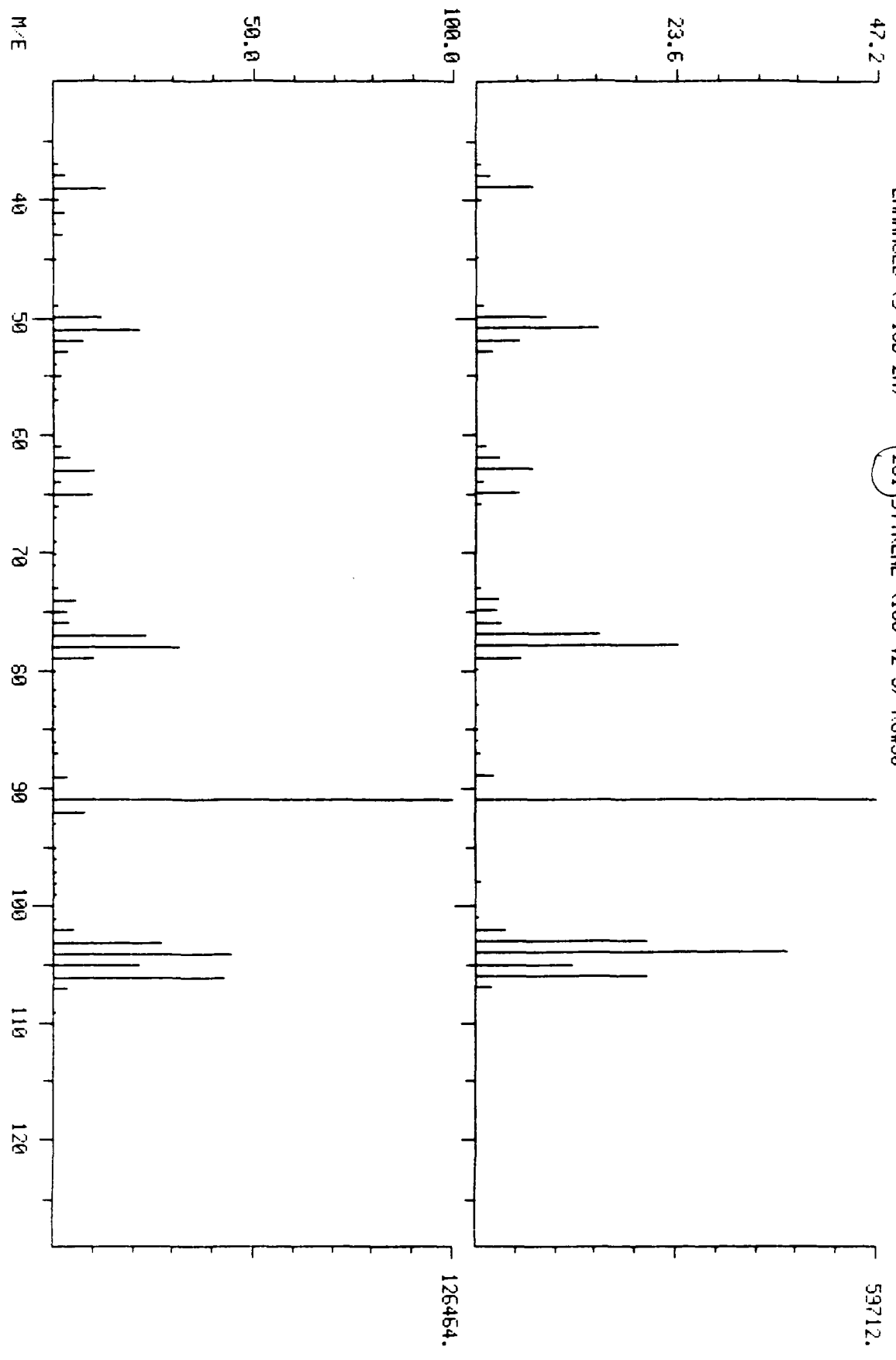
C8.H8
1 WT 1000
B PK 104
RANK 1
IN 50
PUR 631



DUAL MASS SPECTRUM
12/27/89 5:39:00 + 13:51
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (S 158 2N) (251) STYRENE <100-42-5> R0#50

COMPUCHER LABS

DATA: GR009686C19 #1108 BASE M/E: 91 / 91
RIC: 321535. / 563199.



LIBRARY SEARCH
 12/27/89 5:39:00 + 11:02
 SAMPLE: 1G CC#309686 CASE#18756.7 EPA#R2018 ON#19
 ENHANCED (5 158 2H 0T)

COMFUCHEM LABS

DATA: GR009686C19 # 893

BASE M/E: 95
 RIC: 72191.

1000
 SAMPLE

C6.H6.02
 M WT 110
 B PK 95
 RANK 1
 IN 1598
 PUR 835

ETHANONE,1-(2-FURANYL)- CAS# 1192-62-7

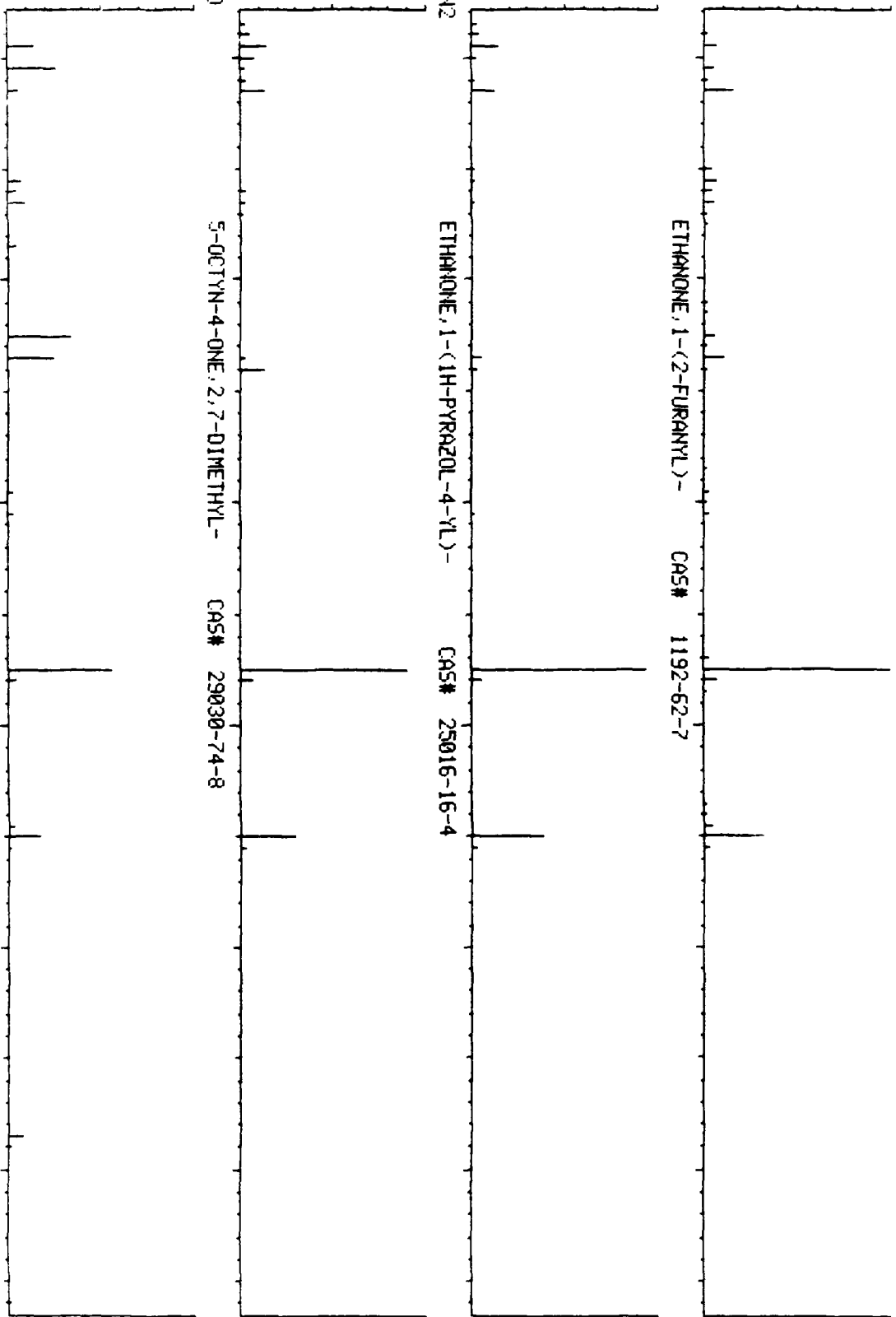
C5.H6.0.N2
 M WT 110
 B PK 95
 RANK 1
 IN 1588
 PUR 787

ETHANONE,1-(1H-PYRAZOL-4-YL)- CAS# 25016-16-4

C10.H16.0
 M WT 190
 B PK 95
 RANK 3
 IN 6245
 PUR 716

5-OCTYN-4-ONE,2,7-DIMETHYL- CAS# 29030-74-8

M/E 40 60 80 100 120 140

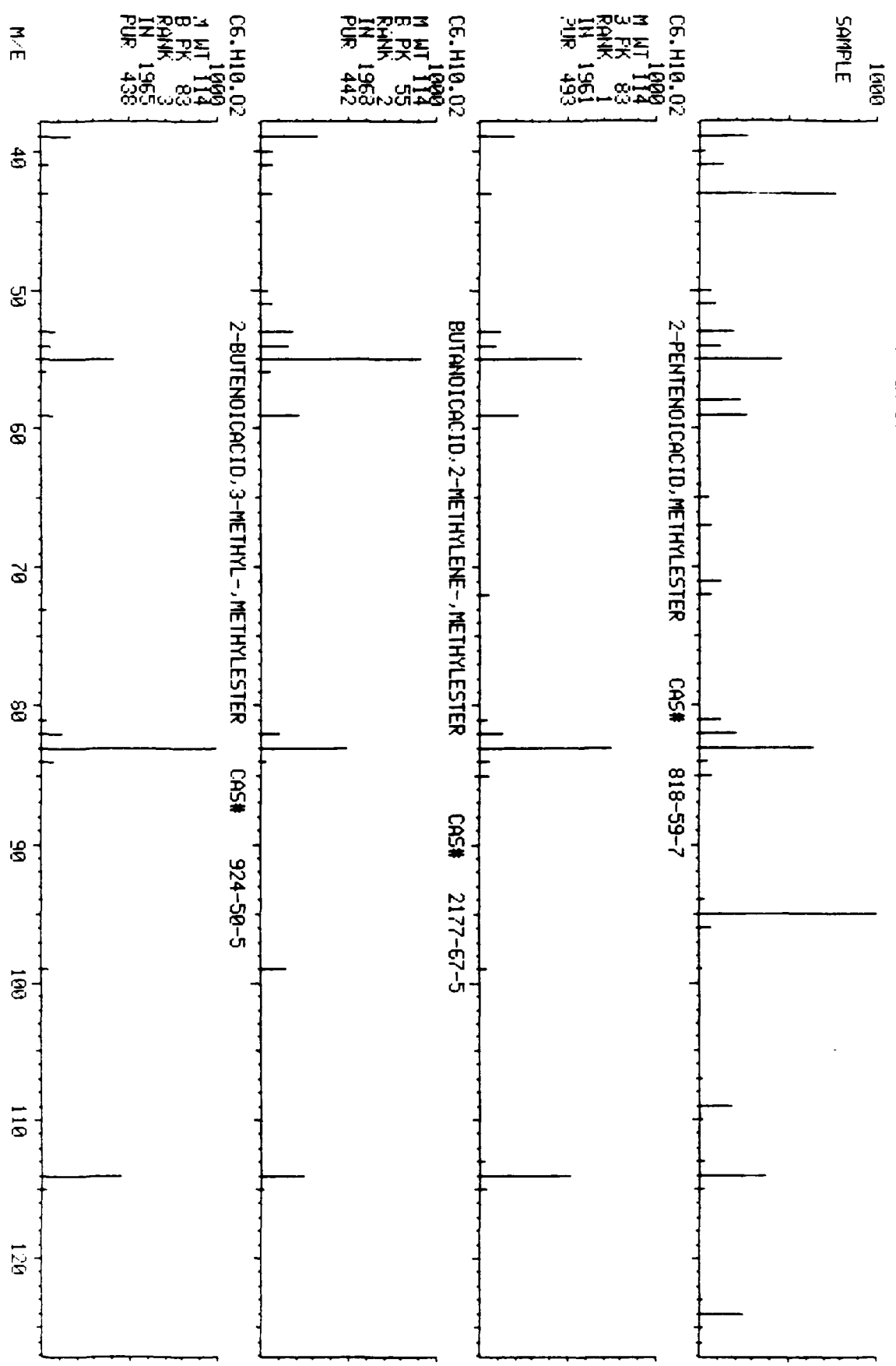


LIBRARY SEARCH
 12/27/89 5:39:00 + 13:27
 SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B2018 ON#19
 ENHANCED (S 158 2N 0T)

COMPUchem LABS

DATA: GR009686C19 #1076

BASE M/E: 95
 RIC: 97535.



LIBRARY SEARCH
12/27/89 5:39:00 + 14:05
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#R2018 ON#19
ENHANCED (5 15B 2N 0T)

COMPUJEN LABS

DATA: GR009686C19 #1127

BASE M/E: 57
RIC: 137727.

1320
SAMPLE

C7.H14.0
M WT 1320
B PK 114
RANK 57
IN 2024
PUR 680

3-HEXANONE,5-METHYL- CAS# 623-56-3

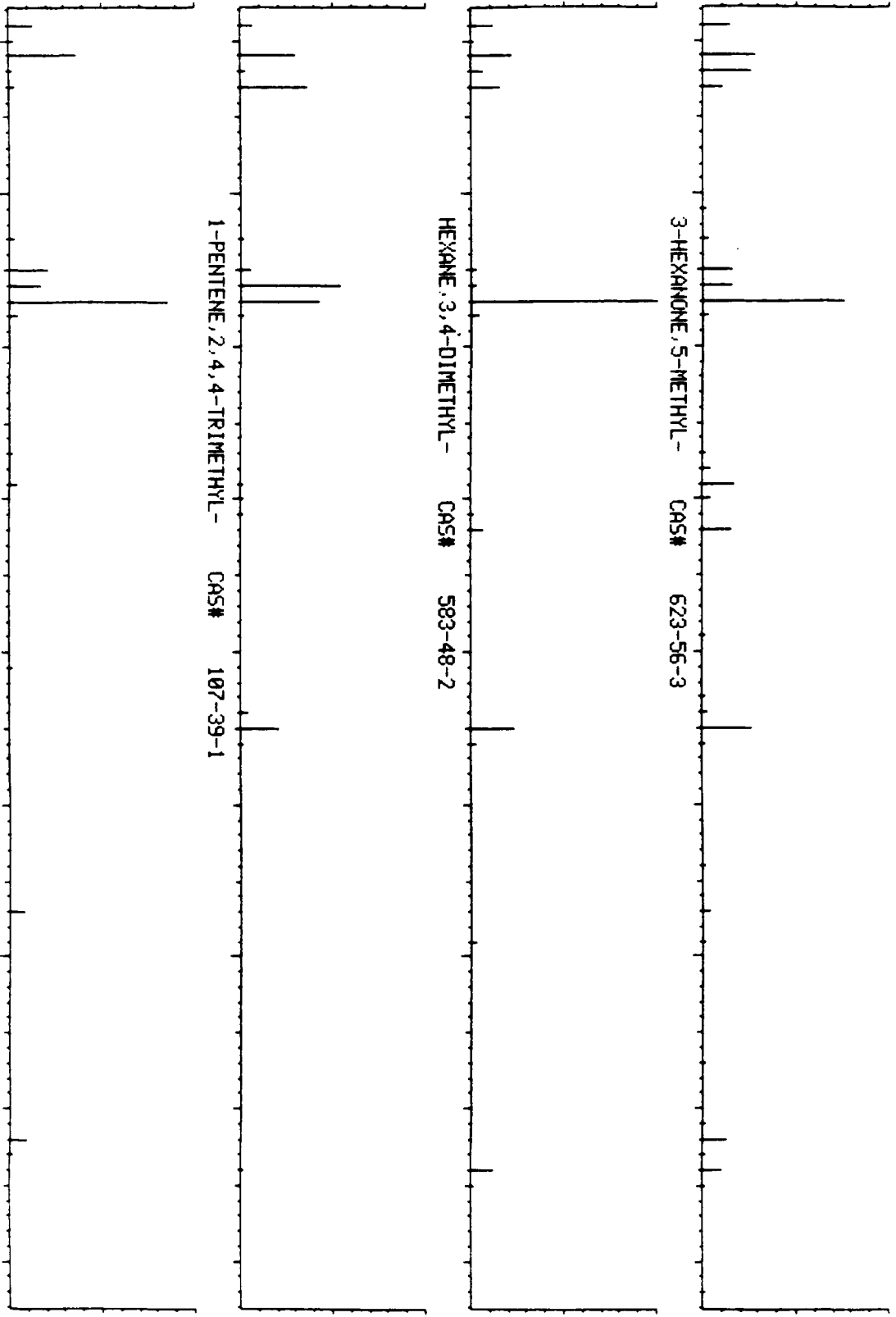
C8.H18
M WT 1320
B PK 114
RANK 56
IN 2079
PUR 606

HEXANE,3,4-DIMETHYL- CAS# 583-48-2

C8.H16
M WT 1320
B PK 112
RANK 57
IN 1814
PUR 606

1-PENTENE,2,4,4-TRIMETHYL- CAS# 107-39-1

M/E 40 50 60 70 80 90 100 110 120



LIBRARY SEARCH
 12/27/89 5:39:00 + 14:16
 SAMPLE: 1G CC#009686 CASE#18756.7 EPA#B201B ON#19
 ENHANCED (5 15B 2H 0T)

COMPUCHEM LABS

DATA: GR009686C19 #1142
 BASE M/E: 43
 RIC: 141823.

1166
 SAMPLE

C7.H14.0
 M WT 1165
 B PK 43
 RANK 1
 IN 2015
 PUR 834

C7.H14.0
 M WT 1165
 B PK 43
 RANK 2
 IN 2014
 PUR 804

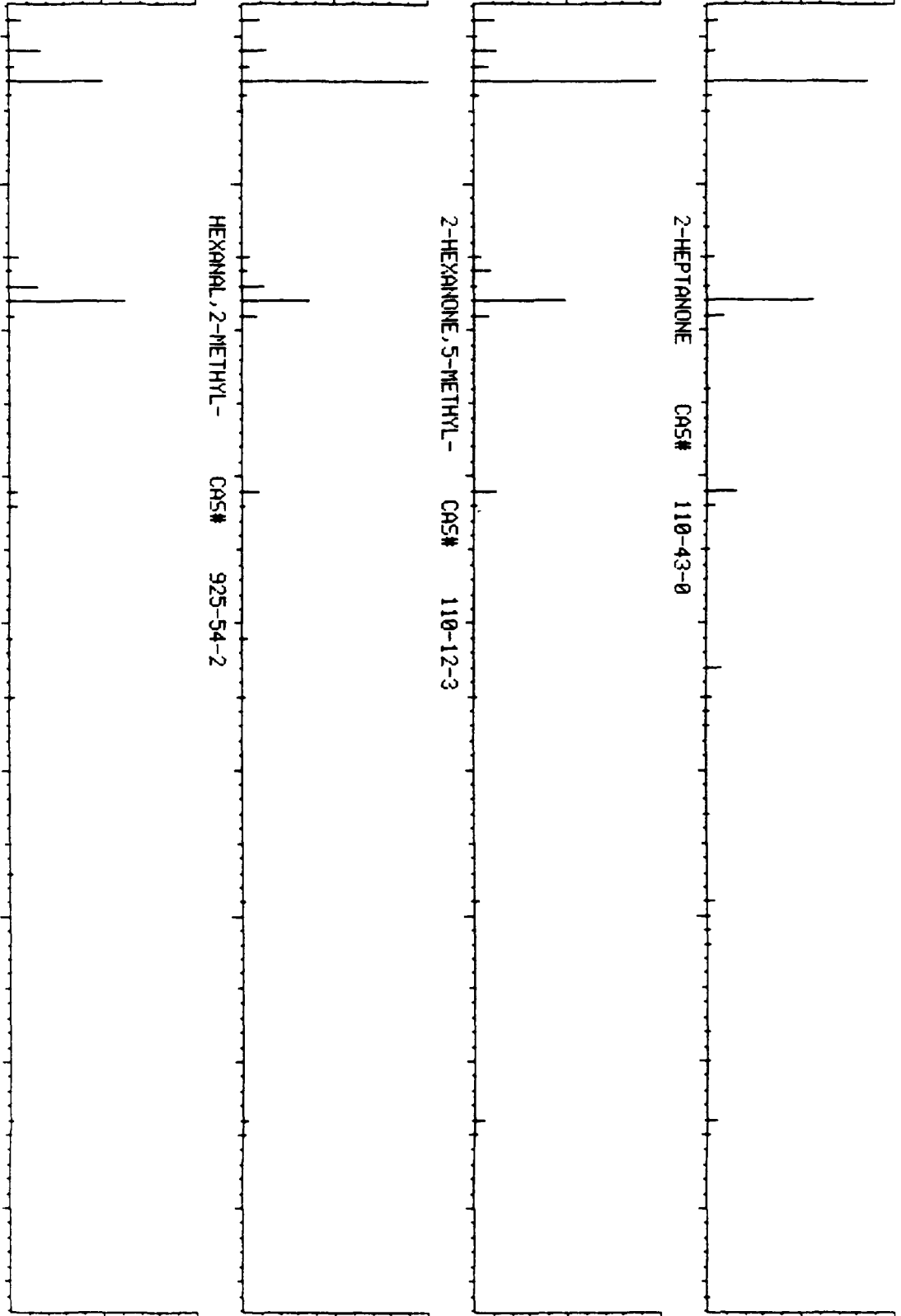
C7.H14.0
 M WT 1165
 B PK 58
 RANK 3
 IN 2025
 PUR 734

M/E

2-HEPTANONE CAS# 110-43-0

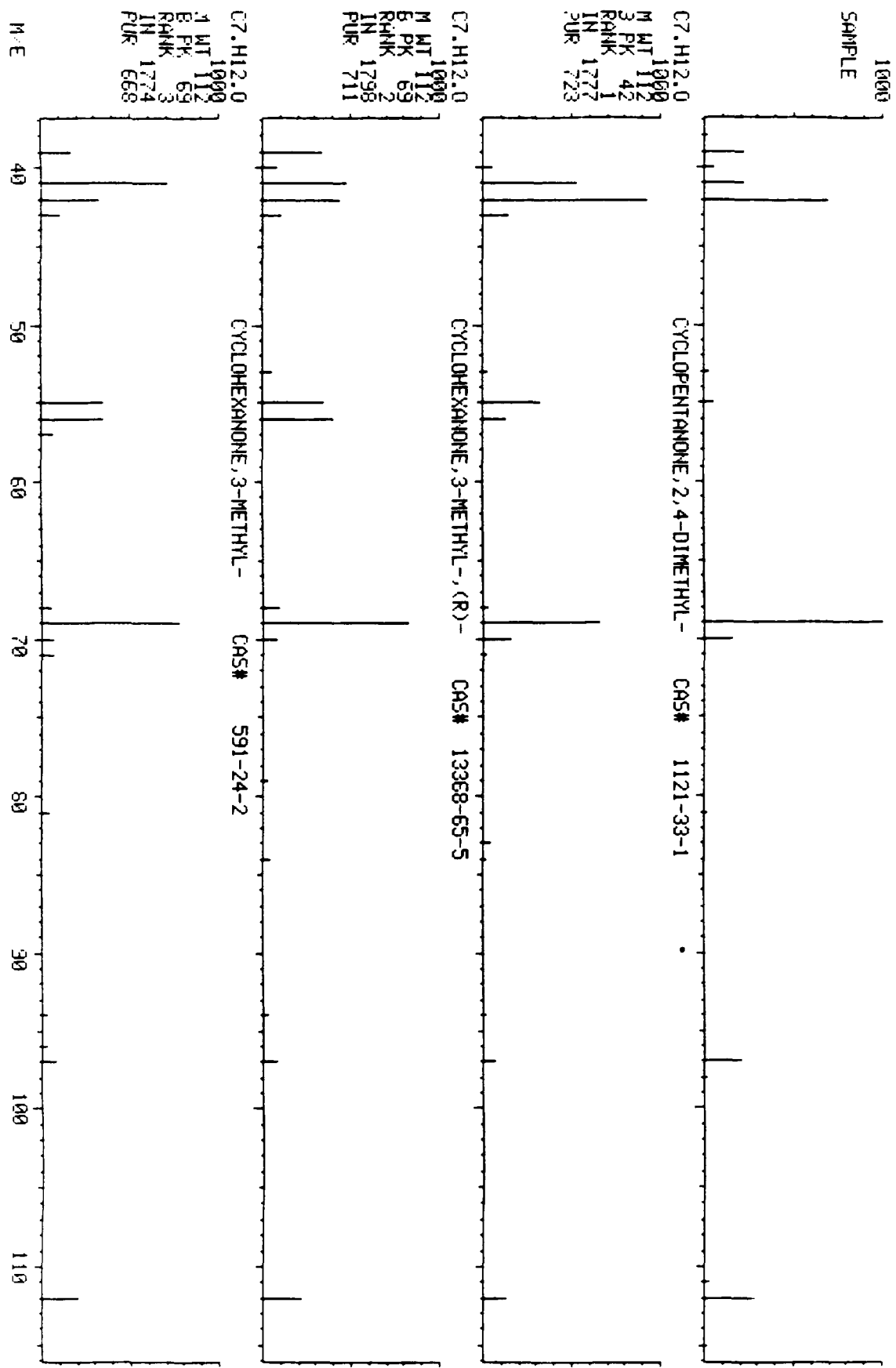
2-HEXANONE,5-METHYL- CAS# 110-12-3

HEXANAL,2-METHYL- CAS# 925-54-2



COMPUCHEM LABS
 LIBRARY SEARCH
 12/27/89 5:39:00 + 14:25
 SAMPLE: 16 CC#309686 CASE#18756.7 EPA#B2018 QM#19
 ENHANCED (5.158 2N 0T)

DATA: GR009686C19 #1154
 BASE N/E: 69
 RIC: 50055.



LIBRARY SEARCH
12/27/89 5:39:00 + 14:33
SAMPLE: 16 GC#309686 CASE#18756.7 EPA#B2018 QN#19
ENHANCED (5 158 21 0T)

COMPUchem LABS

DATA: GR005696013 #1164

BASE M/E: 105
RIC: 51775.

1629
SAMPLE

C7.H12.0

M WT 1629
PK 55
RANK 1
IN 1773
PUR 498

CYCLOHEXANONE, 4-METHYL- CAS# 589-92-4

C7.H12.0

M WT 1629
PK 56
RANK 2
IN 1786
PUR 445

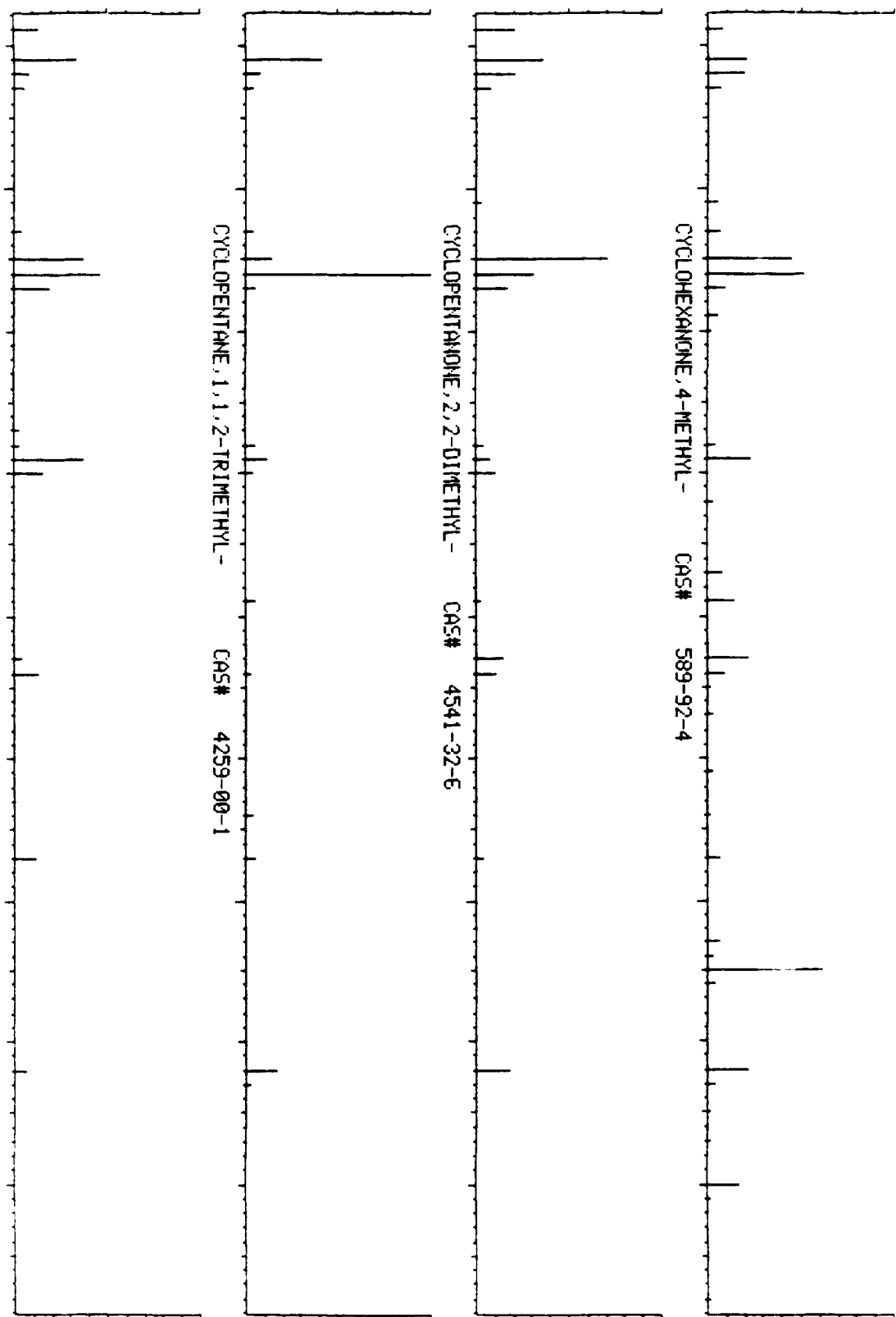
CYCLOPENTANONE, 2,2-DIMETHYL- CAS# 4541-32-6

C8.H16

M WT 1629
PK 56
RANK 3
IN 1844
PUR 440

CYCLOPENTANE, 1,1,2-TRIMETHYL- CAS# 4259-00-1

M/E 40 50 60 70 80 90 100 110 120



LIBRARY SEARCH
 12/27/89 5:39:00 + 14:43
 SAMPLE: 16 CC#309586 CASE#18756.7 EPA#B201B ON#19
 ENHANCED (5 156 2N 0T)

COMPUCHEN LHB5

DATA: GR009686C19 #1176

BASE M/E: 108
 RIC: 267311.

1263
 SAMPLE

C7.H8.0

M WT 1253
 B PK 108
 RANK 108
 IN 1505
 PUR 886

BENZENE, METHOXY-

CAS# 100-66-3

C6.H8.N2

M WT 1253
 B PK 108
 RANK 78
 IN 1482
 PUR 745

HYDRAZINE, PHENYL-

CAS# 100-63-0

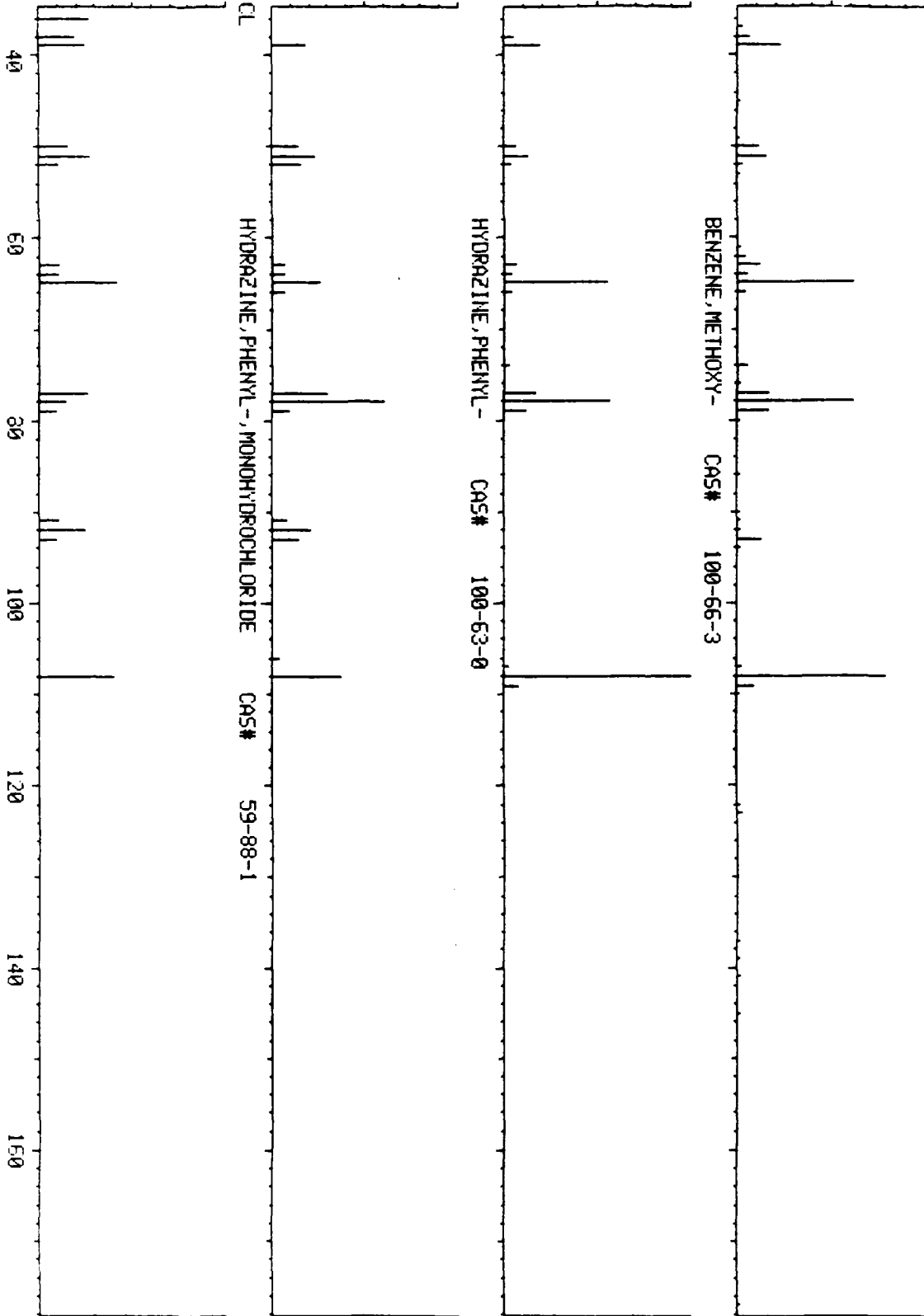
C6.H9.N2.Cl

M WT 1253
 B PK 144
 RANK 65
 IN 5153
 PUR 683

HYDRAZINE, PHENYL-, MONOHYDROCHLORIDE

CAS# 59-88-1

M/E



LIBRARY SEARCH
 12/27/89 5:59:00 + 15:21
 SAMPLE: 16 CG#809686 CASE#18756.7 EPA#B2018 QN#19
 ENHANCED (S 158 ZN 01)

COMPUchem LABS

DATA: GR009686C19 #1228

BASE N/E: 91
 RIC: 101119.

1000
 SAMPLE

C16.H14.O3
 1000
 M WT 254
 R PK 91
 RANK 1
 IN 18395
 PUR 675

C8.H8.O
 1000
 M WT 128
 R PK 91
 RANK 2
 IN 2517
 PUR 630

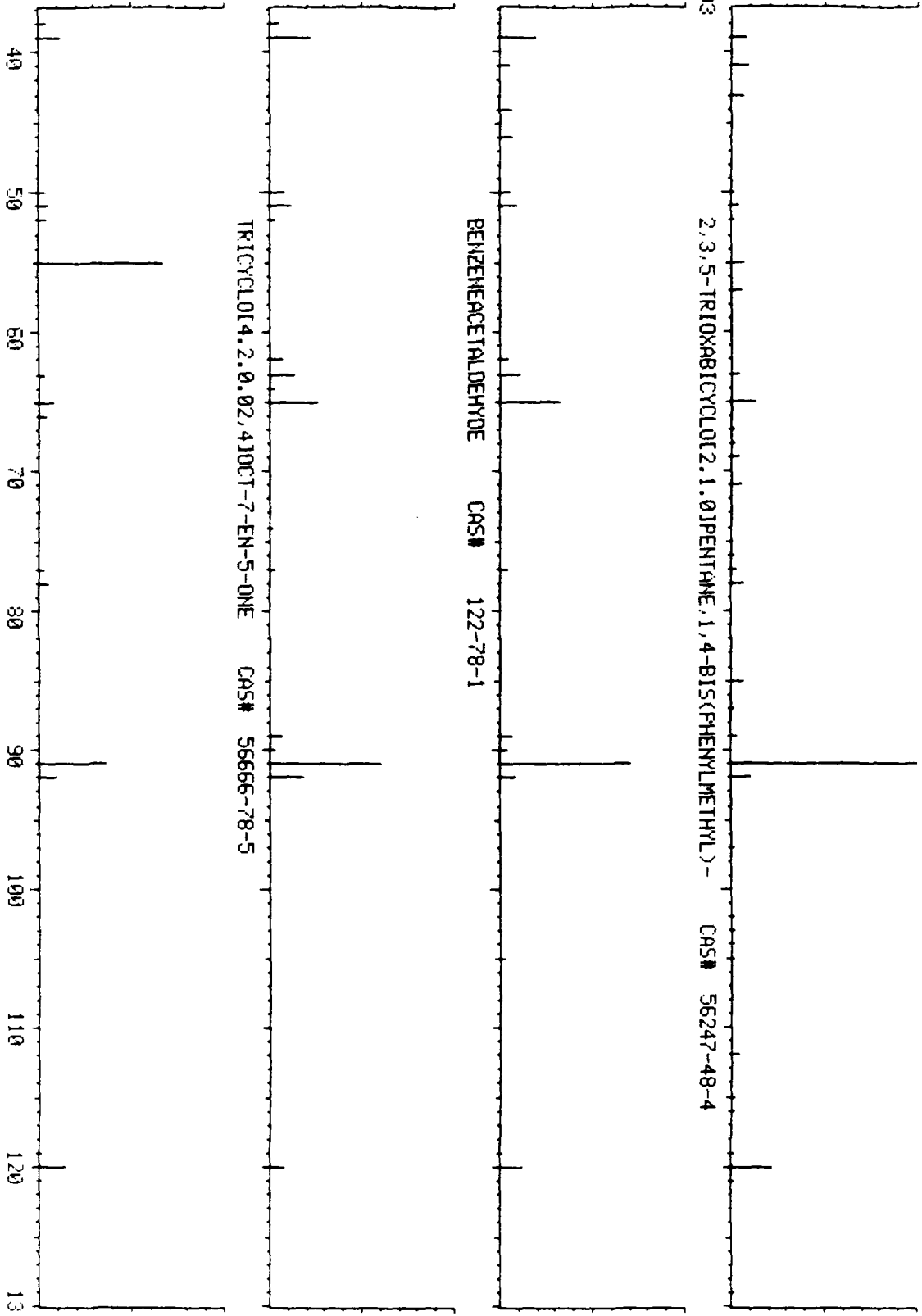
8.H8.O
 1000
 M WT 128
 R PK 55
 RANK 3
 IN 2522
 PUR 600

M/E

2,3,5-TRIOXABICYCLO[2.1.0]PENTANE,1,4-BIS(PHENYLMETHYL)- CAS# 56247-48-4

BENZENEACETALDEHYDE CAS# 122-78-1

TRICYCLO[4.2.0.0,2,4]OCT-7-EN-5-ONE CAS# 56666-78-5

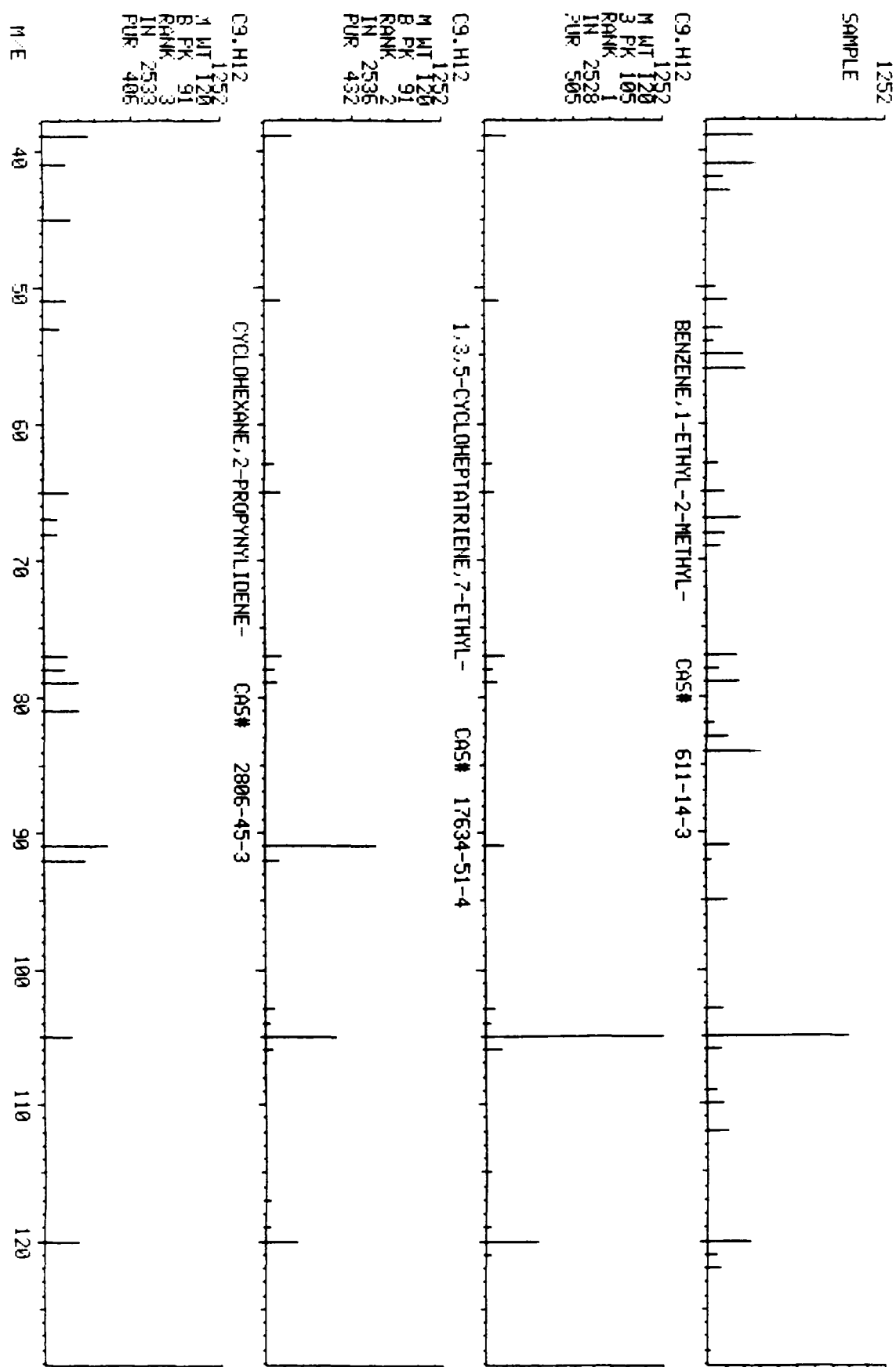


LIBRARY SEARCH
12/27/89 5:39:00 + 15:37
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#R2018 ON#19
ENHANCED (5.156 2N 0T)

COMPUchem LIBS

DATA: GR009686C19 #1249

BASE M/E: 105
RIC: 338943.



LIBRARY SEARCH
12/27/89 5:39:00 + 16:04
SAMPLE: 1G CC#309686 CASE#18756.7 EPA#B201B ON#19
ENHANCED (S 158 2N 0T)

COMPUCHEN LABS
DATA: GR009686C19 #1286

BASE M/E: 105
FIC: 204799.

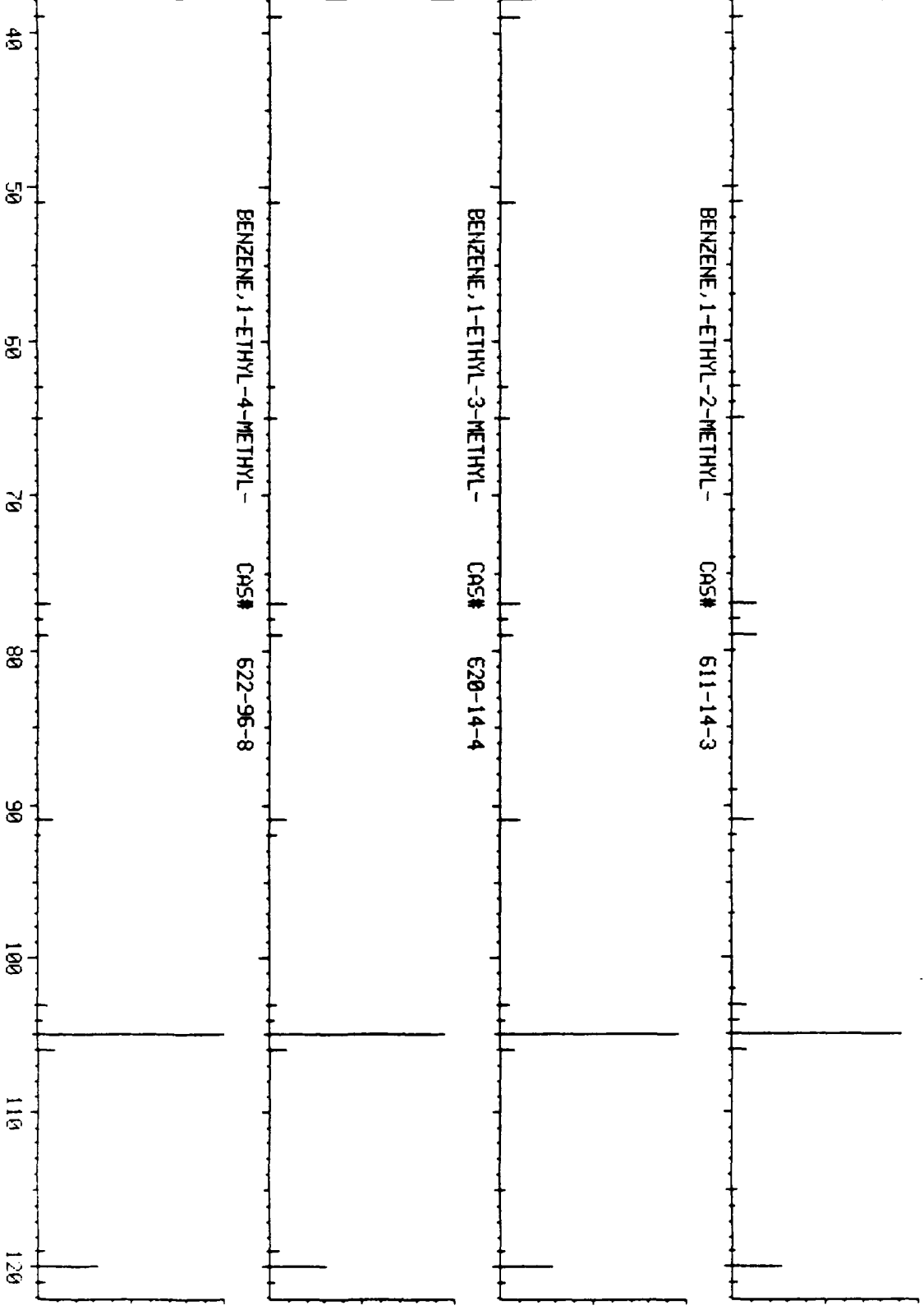
1102
SAMPLE

C9.H12
M WT 1102
3 PK 105
RANK 1
IN 2528
PUR 921

C9.H12
M WT 1102
B PK 105
RANK 2
IN 2529
PUR 912

C9.H12
M WT 1102
B PK 105
RANK 3
IN 2530
PUR 906

M/E



LAB INSTRUCTIONS

SAMPLED DATE 12/18/89

RECEIPT DATE 12/20/89

CASE# 18756 7

DUE DATE

VOA
GC/MS WORKSHEET

COMPUCHEM# 309686R

R[X] R2[] DE []
R3[] R4[] D2[]

L L SOLID, EPA 90W 2/88

Sample Prep Code---155
Instrument Code---413
Compound List-----494
Surrogate Std-----394
Internal Std-----038

=====

SAMPLE ID#: B201B Dry Wt. Factor 1.14 % Moisture 1

=====

GC/MS ANALYSIS

Amount Purged: [/] 10mls/Xg soil or [] Dilution _____ ul/10000ul/Xg soil
Internal Standard Volume Added 5 ul
Surrogate Standard Volume Added 5 ul
BFB Filename RE891227/19 Disk (2958A)
Blank Filename H010660/19 Disk ()
Standard Filename 61891227/19 Disk ()
Sample Filename CR009486/19 Disk ()

ANALYST(S): Injection 1477-16 Work-up 1477-16

GC/MS REVIEW

CONDITION
CODE

DI

Entry Codes OK, EA, ES, SM, JS, SL, SH, JA, DA

Non-Entry Codes IM, IL, IH, SW, CT, CS, PC, NP
IF, LA, DI, CO, RN, DW, SI, SF
UP, BB, OT, VC, FO, NS

Disposition: [/] Complete
[] Reprep neat required
[] Reprep using _____
[] Dilute ()
1/5 Reprep Med.

Extraneous Peak Search Results:

of Peaks Found: 10

MM
1/30/90

Quality Assurance Notice(s):

Notices Required 0

COMMENTS:

GC/MS Review EDWagner Date 1/2/90 Auditor EDWagner Date 1/3/90

REPORT INTEGRATION

Final Reportable Package(s) GR0-19

Total # of Injections 6

1 CSR-13

QA COMMENTS

FINAL REVIEW

Initials _____ Date _____

Initials _____ Date _____

AC1007 105 BT

VOLATILE PREPARATION WORKSHEET

QUEUE # 19 PREP CODE 155 ASSIGNED TO Harshad Jeshi DATE 12/22/89

SAMPLE NUMBER	CASE NUMBER	SDG	QC SAMPLE		SAMPLE WEIGHT (G) VOLUME (ML)	SAMPLE ID	COMMENTS
			TYPE	ORIGINAL			
309679R	18756	0007			1.0g	B201A	
309680R			SS	309679	1.0g	B201A MS	
309681R			SS	309679	1.0g	B201A MSD	
309682R			BS		0.0g		
309686R					1.0g	B201B	
309687R					1.0g	B202TAR	
310543	18244	0142			5.0g	MW-17-30	
310555	18477	0132			5.0g	EPTP10122	
310560			SS	310555	5.0g	EPTP10122 MS	
310561			SS	310555	5.0g	EPTP10122 MSD	
310562			BS		0.0g		
310659			B1		5.0ml	F4	
310660			B2		0.0ml	F5	
310661			B3		0.0ml	F6	
310662			B4		0.0ml	F7	
310663			B5		0.0ml	F8	

SURROGATE # / LOT # / MANUAL OPERATOR F3, 1518
 AMOUNT
 RELINQUISHED BY Indo DATE 12/21/89 RECEIVED BY 182 Naptano DATE 12/21/89

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	464	52500	50.0		X57
221	50	CHLOROMETHANE				BDL	57 60-9
231	62	VINYL CHLORIDE				BDL	5
220	94	BROMOMETHANE				BDL	5
209	64	CHLOROETHANE				BDL	5
216	96	1,1-DICHLOROETHENE				BDL	28 4-13
254	76	CARBON DISULFIDE				BDL	5
252	43	ACETONE (2-PROPANONE)			64.8	320 320 E	5
248	114 I	1,4-DIFLUOROBENZENE (IS)	610	198000	50.0		
222	84	METHYLENE CHLORIDE			12.1	60 65 B	20
226	96	TRANS-1,2-DICHLOROETHENE				BDL	20
214	63	1,1-DICHLOROETHANE				BDL	20
257	43	VINYL ACETATE				BDL	20
237	96	CIS-1,2-DICHLOROETHENE				BDL	20
253	72	2-BUTANONE			46.4	230 260	20
211	83	CHLOROFORM				BDL	20
227	97	1,1,1-TRICHLOROETHANE				BDL	20
206	117	CARBON TETRACHLORIDE				BDL	20
203	78	BENZENE			25.3	130 140	20
215	62	1,2-DICHLOROETHANE				BDL	20
270	117 I	D5-CHLOROBENZENE (IS) RO#29	1001	195000	50.0		
229	130	TRICHLOROETHENE				BDL	20
217	63	1,2-DICHLOROPROPANE			1.3	BDL 65	20
212	83	BROMODICHLOROMETHANE				BDL	20
218	75	CIS-1,3-DICHLOROPROPENE				BDL	20
256	43	4-METHYL-2-PENTANONE			12.5	62 71	5
225	92	TOLUENE			280.0	1400 1400 E	20
250	75	TRANS-1,3-DICHLOROPROPENE				BDL	20
228	97	1,1,2-TRICHLOROETHANE				BDL	20
224	164	TETRACHLOROETHENE				BDL	20
255	43	2-HEXANONE			273.0	1400 1400 E B	50
208	129	DIBROMOCHLOROMETHANE				BDL	20
207	112	CHLOROBENZENE				BDL	20
219	106	ETHYLBENZENE			172.0	860 980	20
330	106	M, P-XYLENE			337.0	1700 1700 E	20
239	106	O-XYLENE			249.0	1200 1200 E	20
251	104	STYRENE			95.3	480 480	20
205	173	BROMOFORM				BDL	20
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	20
258	65 S	D4-1,2-DICHLOROETHANE RO#57			47.2	94. %	
247	95 S	BROMOFLUOROBENZENE			49.7	99. %	
233	98 S	D8-TOLUENE RO#59			44.9	90. %	
289	106	XYLENES (TOTAL)			587.0	2900 2900 E	20

CORRECTED/REVIEWED BY *OK*
(GC/MS DATA REVIEWER)

DATE 1-2-90

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
299	96	1,2-DICHLOROETHENE (TOTAL)				BDL	2
CHECKSUMS:							
	0979		2075	445500.	2447.5	11031	

CORRECTED/REVIEWED BY

OK [Signature]
(GC/MS DATA REVIEWER)

DATE

1.29.90

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P
40	258	D4-1,2-DICHLOROETHANE RO#57	47.2	50.0	94.	70-121	X
41	247	BROMOFLUOROBENZENE	49.7	50.0	99.	74-121	X
42	233	D8-TOLUENE RO#59	44.9	50.0	90.	81-117	X

* ADVISORY SURROGATE ONLY

++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

=====

CORRECTION FACTOR CALCULATION:

$$\frac{5.0 \text{ G}}{\text{WET WEIGHT OF SAMPLE (G)}} \times \frac{\text{GC/MS DILUTION FACTOR}}{\text{GC/MS DILUTION FACTOR}} \times \text{DRY WEIGHT FACTOR} =$$

$$\frac{5.0 \text{ G}}{1.00 \text{ (G)}} \times \frac{1.0}{1.00} \times \frac{1.14}{1.00} = \frac{5.76}{5.000}$$

=====

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:

THE SURROGATES ARE ADDED TO THE SAMPLE PRIOR TO SPARGING.
SURROGATE SPIKE CONVERSION FACTOR = 1.

=====

VERSION 8

CORRECTED/REVIEWED BY *C. J. + S.*
(GC/MS DATA REVIEWER)

DATE 1-2-90

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B201BRE

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C5R09686B13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/29/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	1400	U
74-83-9	-----Bromomethane	1400	U
75-01-4	-----Vinyl Chloride	1400	U
75-00-3	-----Chloroethane	1400	U
75-09-2	-----Methylene Chloride	540	DBJ
67-64-1	-----Acetone	1400	U
75-15-0	-----Carbon Disulfide	710	U
75-35-4	-----1,1-Dichloroethene	710	U
75-34-3	-----1,1-Dichloroethane	710	U
540-59-0	-----1,2-Dichloroethene (total)	710	U
67-66-3	-----Chloroform	710	U
107-06-2	-----1,2-Dichloroethane	710	U
78-93-3	-----2-Butanone	1400	U
71-55-6	-----1,1,1-Trichloroethane	710	U
56-23-5	-----Carbon Tetrachloride	710	U
108-05-4	-----Vinyl Acetate	1400	U
75-27-4	-----Bromodichloromethane	710	U
78-87-5	-----1,2-Dichloropropane	710	U
10061-01-5	-----cis-1,3-Dichloropropene	710	U
79-01-6	-----Trichloroethene	710	U
124-48-1	-----Dibromochloromethane	710	U
79-00-5	-----1,1,2-Trichloroethane	710	U
71-43-2	-----Benzene	2200	D
10061-02-6	-----Trans-1,3-Dichloropropene	710	U
75-25-2	-----Bromoform	710	U
108-10-1	-----4-Methyl-2-Pentanone	1400	U
591-78-6	-----2-Hexanone	2400	D
127-18-4	-----Tetrachloroethene	710	U
79-34-5	-----1,1,2,2-Tetrachloroethane	710	U
108-88-3	-----Toluene	16000	D
108-90-7	-----Chlorobenzene	710	U
100-41-4	-----Ethylbenzene	7400	D
100-42-5	-----Styrene	2800	D
1330-20-7	-----Total Xylenes	21000	D

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

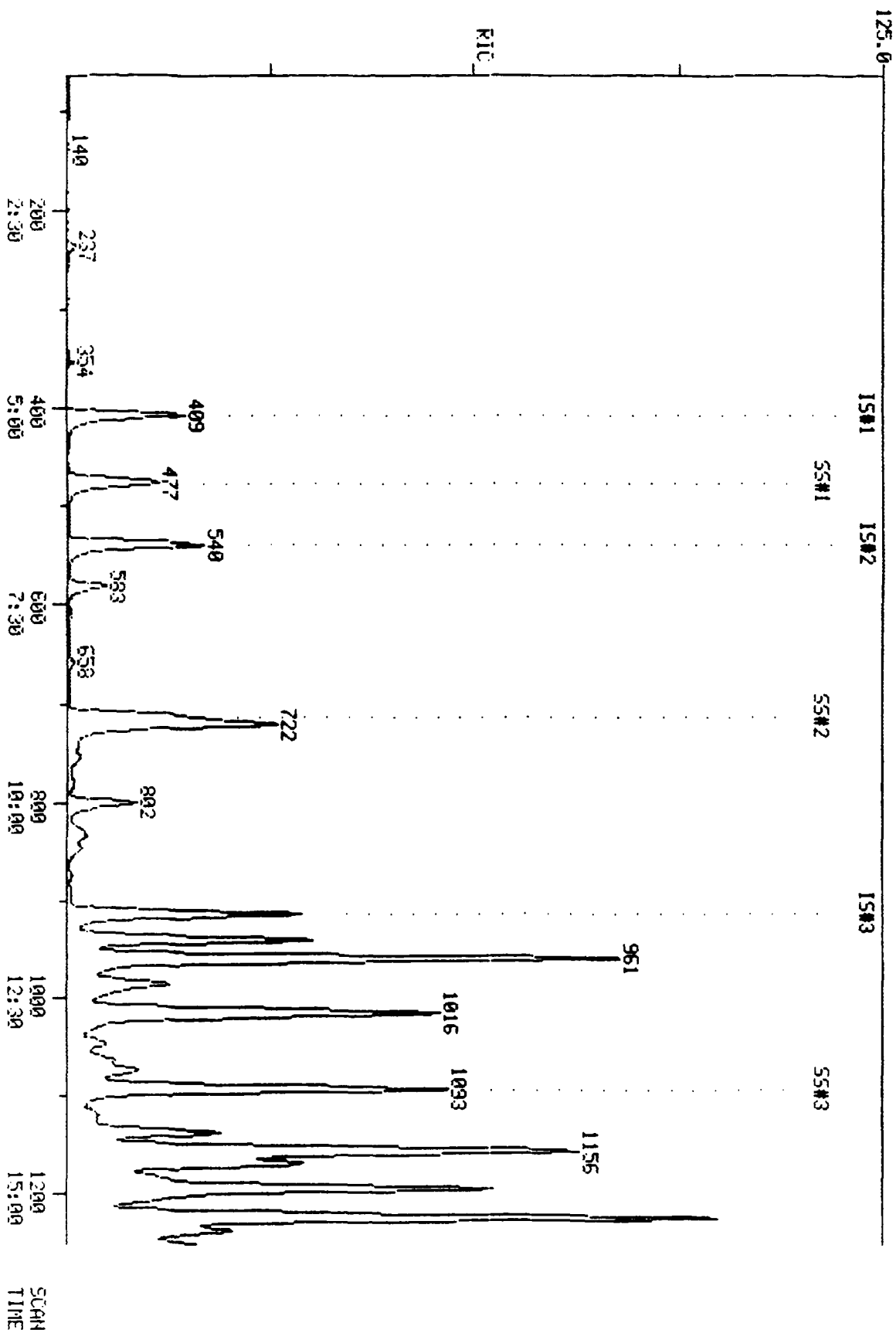
B201BRE

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309686
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C5R09686B13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/29/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 625-86-5	FURAN, 2,5-DIMETHYL-	7.28	1800	J
2.	UNKNOWN	10.02	2000	J
3.	UNKNOWN	12.34	2800	J
4.	METHYLETHYLBENZENE + UNKNOWN	13.44	710	J
5. 103-65-1	BENZENE, PROPYL-	14.22	3100	J
6. 98-82-8	BENZENE, (1-METHYLETHYL)-	14.45	17000	J
7.	TRIMETHYLBENZENE ISOMER	14.60	4700	J
8.	ETHYLMETHYLBENZENE + UNKNOWN	14.94	14000	J
9.	TRIMETHYLBENZENE ISOMER	15.30	18000	J
10. 611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	15.45	2000	J

RLC
 12/29/89 20:10:00
 SAMPLE: 100UL CC#309686 EPAN#Z01B CASE#18756 ON #13
 COND5.:
 COMPUTHER LABS
 RE: *CS# 11210*
 COMPUTHER DATA: CS#09686R13 SCANS 65 TO 1250
 359040.



QUANTITATION REPORT FILE: C5R09686B13
 DATA: C5R09686B13.T1
 12/29/89 20:10:00
 SAMPLE: 100UL CC#309686 EPA#B201B CASE#18756 ON #13
 CONDS: 54 sub, 1/2/90
 SUBMITTED BY: 13 ANALYST: 13

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
 RESP. FAC. FROM LIBRARY ENTRY

- | NO | NAME |
|----|--|
| 1 | *234 BROMOCHLOROMETHANE (IS) <75-97-5> NA#1 |
| 2 | 221 CHLOROMETHANE <74-87-3> NA#2 |
| 3 | 231 VINYL CHLORIDE <75-01-4> NA#3 |
| 4 | 220 BROMOMETHANE <78-83-9> NA#4 |
| 5 | 209 CHLOROETHANE <75-00-3> NA#5 |
| 6 | 216 1,1-DICHLOROETHENE <75-35-4> NA#8 |
| 7 | 254 CARBON DISULFIDE <75-15-0> NA#9 |
| 8 | 252 ACETONE (2-PROPANONE) <67-64-1> NA#13 |
| 9 | *248 1,4-DIFLUOROBENZENE (IS) <540-36-3> NA#14 |
| 10 | 222 METHYLENE CHLORIDE <75-09-2> NA#16 |
| 11 | 226 TRANS-1,2-DICHLOROETHENE <156-60-5> NA#17 |
| 12 | 214 1,1-DICHLOROETHANE <75-34-3> NA#19 |
| 13 | 257 VINYL ACETATE <108-05-4> NA#20 |
| 14 | 237 CIS-1,2-DICHLOROETHENE <156-59-2> NA#21 |
| 15 | 253 2-BUTANONE <78-93-3> NA#22 |
| 16 | 211 CHLOROFORM <67-66-2> NA#23 |
| 17 | 227 1,1,1-TRICHLOROETHANE <71-55-6> NA#24 |
| 18 | 206 CARBON TETRACHLORIDE <56-23-5> NA#25 |
| 19 | 203 BENZENE <71-43-2> NA#26 |
| 20 | 215 1,2-DICHLOROETHANE <107-06-2> NA#27 |
| 21 | *270 D5-CHLOROENZENE (IS) <XX-XX-X> NA#29 |
| 22 | 229 TRICHLOROETHENE <79-01-6> NA#30 |
| 23 | 217 1,2-DICHLOROPROPANE <78-87-5> NA#31 |
| 24 | 212 BROMODICHLOROMETHANE <75-27-4> NA#33 |
| 25 | 218 CIS-1,3-DICHLOROPROPENE <10061-1-5> NA#35 |
| 26 | 256 4-METHYL-2-PENTANONE <108-01-1> NA#36 |
| 27 | 225 TOLUENE <108-88-3> NA#37 |
| 28 | 250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> NA#38 |
| 29 | 228 1,1,2-TRICHLOROETHANE <79-00-5> NA#39 |
| 30 | 224 TETRACHLOROETHENE <127-18-4> NA#41 |
| 31 | 255 2-HEXANONE <391-78-6> NA#42 |
| 32 | 208 DIBROMOCHLOROMETHANE <124-48-1> NA#43 |
| 33 | 207 CHLOROBENZENE <108-90-7> NA#45 |
| 34 | 219 ETHYLBENZENE <100-41-4> NA#47 |
| 35 | 330 M,P-XYLENE <133-02-7> NA#48 |
| 36 | 239 O-XYLENE <133-02-7> NA#49 |
| 37 | 251 STYRENE <100-42-5> NA#50 |
| 38 | 205 BROMOFORM <75-25-2> NA#51 |
| 39 | 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> NA#54 |
| 40 | *258 D4-1,2-DICHLOROETHANE NA#57 |
| 41 | *247 BROMOFLUOROBENZENE <460-00-4> NA#58 |
| 42 | *233 D8-TOLUENE NA#59 |

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	409	5.07	1	1.000	A BB	54485.	50.000 UG/L	7 29
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	199	2:29	1	0.487	A BV	367.	1.436 UG/L	0.21
9	114	540	6:45	9	1.000	A BB	189860.	50.000 UG/L	7.29
10	84	238	2:58	1	0.582	A BB	5930.	3.790 UG/L	0.55
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	NOT FOUND							
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	474	5:55	9	0.878	A BB	45195.	15.249 UG/L	2.22
20	62	NOT FOUND							
21	117	914	11:25	21	1.000	A BB	269834.	50.000 UG/L	7.29
22	130	NOT FOUND							
23	63	NOT FOUND							
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	724	9:03	21	0.792	A VB	2277.	2.042 UG/L	0.30
27	92	723	9:02	21	0.791	A BB	156888.	111.890 UG/L	16.31
28	75	NOT FOUND							
29	97	NOT FOUND							
30	164	NOT FOUND							
31	43	847	10:35	21	0.927	A VB	18130.	16.819 UG/L	2.45
32	129	NOT FOUND							
33	112	NOT FOUND							
34	106	942	11:46	21	1.031	A BV	105991.	52.018 UG/L	7.58
35	106	960	12:00	21	1.050	A VB	310656.	86.631 UG/L	12.63
36	106	1015	12:41	21	1.111	A BB	189103.	60.376 UG/L	8.80
37	104	1021	12:46	21	1.117	A BB	105821.	19.642 UG/L	2.86
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	477	5:58	1	1.166	A BB	110676.	56.352 UG/L	8.22
41	95	1094	13:40	21	1.197	A BB	247443.	52.302 UG/L	7.63
42	98	714	8:55	21	0.781	A BB	145509.	57.294 UG/L	8.35

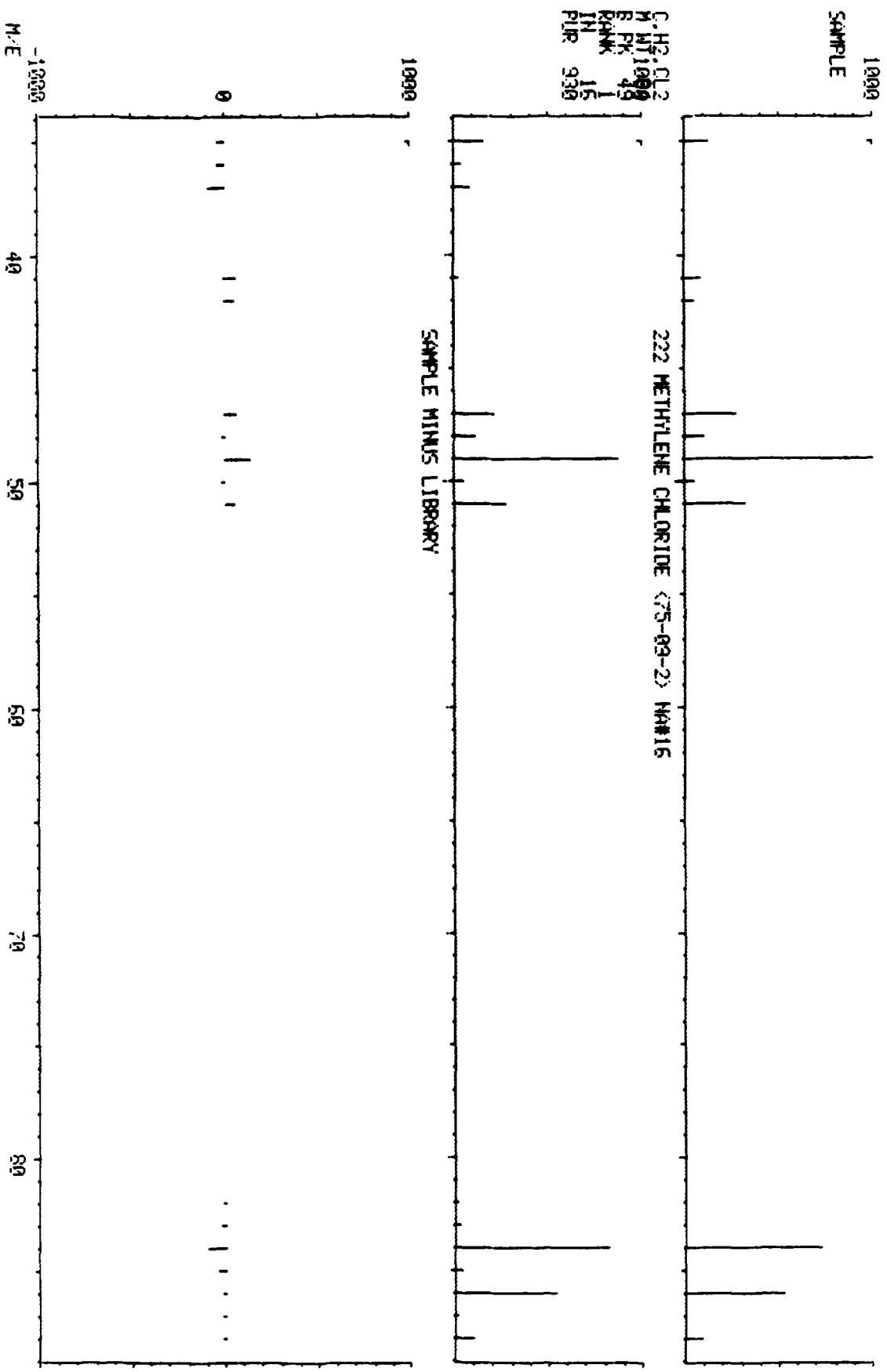
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:08	1.00	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	0:50		10.000			50.00		0.551	
3	0:57		10.000			50.00		0.697	
4	1:11		10.000			50.00		0.985	
5	1:17		10.000			50.00		0.487	
6	2:10		5.000			50.00		1.108	
7	2:17		5.000			50.00		3.396	
8	2:31	0.99	10.000	0.05	1.44	50.00	0.007	0.235	0.03
9	6:49	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	2:58	1.00	5.000	0.12	3.79	50.00	0.109	1.436	0.08
11	3:19		5.000			50.00		1.159	
12	3:56		5.000			50.00		1.953	
13	4:16		10.000			50.00		0.803	
14	4:49		5.000			50.00		1.474	

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:04		10.000			50.00		0.137	
16	5:24		5.000			50.00		2.637	
17	5:25		5.000			50.00		0.654	
18	5:37		5.000			50.00		0.592	
19	5:59	0.99	5.000	0.18	15.25	50.00	0.238	0.781	0.30
20	6:08		5.000			50.00		1.645	
21	11:32	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	7:04		5.000			50.00		0.387	
23	7:25		5.000			50.00		0.311	
24	7:59		5.000			50.00		0.547	
25	8:43		5.000			50.00		0.467	
26	9:10	0.99	15.000	0.05	2.04	50.00	0.008	0.207	0.04
27	9:07	0.99	5.000	0.16	111.89	50.00	0.581	0.260	2.24
28	9:49		5.000			50.00		0.369	
29	10:04		5.000			50.00		0.516	
30	10:04		5.000			50.00		0.512	
31	10:40	0.99	15.000	0.06	16.82	50.00	0.067	0.200	0.34
32	10:38		5.000			50.00		0.809	
33	11:35		5.000			50.00		0.867	
34	11:53	0.99	5.000	0.21	52.02	50.00	0.393	0.378	1.04
35	12:07	0.99	5.000	0.21	86.63	50.00	1.151	0.664	1.73
36	12:48	0.99	5.000	0.22	60.38	50.00	0.701	0.580	1.21
37	12:52	0.99	5.000	0.22	19.64	50.00	0.392	0.998	0.39
38	13:08		5.000			50.00		0.696	
39	14:21		5.000			50.00		0.587	
40	6:01	0.99	5.000	0.23	56.35	50.00	2.031	1.802	1.13
41	13:47	0.99	5.000	0.24	52.30	50.00	0.917	0.877	1.05
42	9:01	0.99	5.000	0.16	57.29	50.00	0.539	0.471	1.15

LIBRARY SEARCH
 12/23/89 20:10:00 + 2:58
 SAMPLE: 100UL CC#309585 EPA#82018 CASE#18765 ON #13
 ENHANCED (S 158 2N 0T)

COMPUchem LABS
 DATA: CSR09686813 # 238
 BASE M/E: 49
 RIC: 3043.

C.H2:CL2
 M.Wt: 100.0
 B.Pk: 49
 IN: 15
 PUP: 930



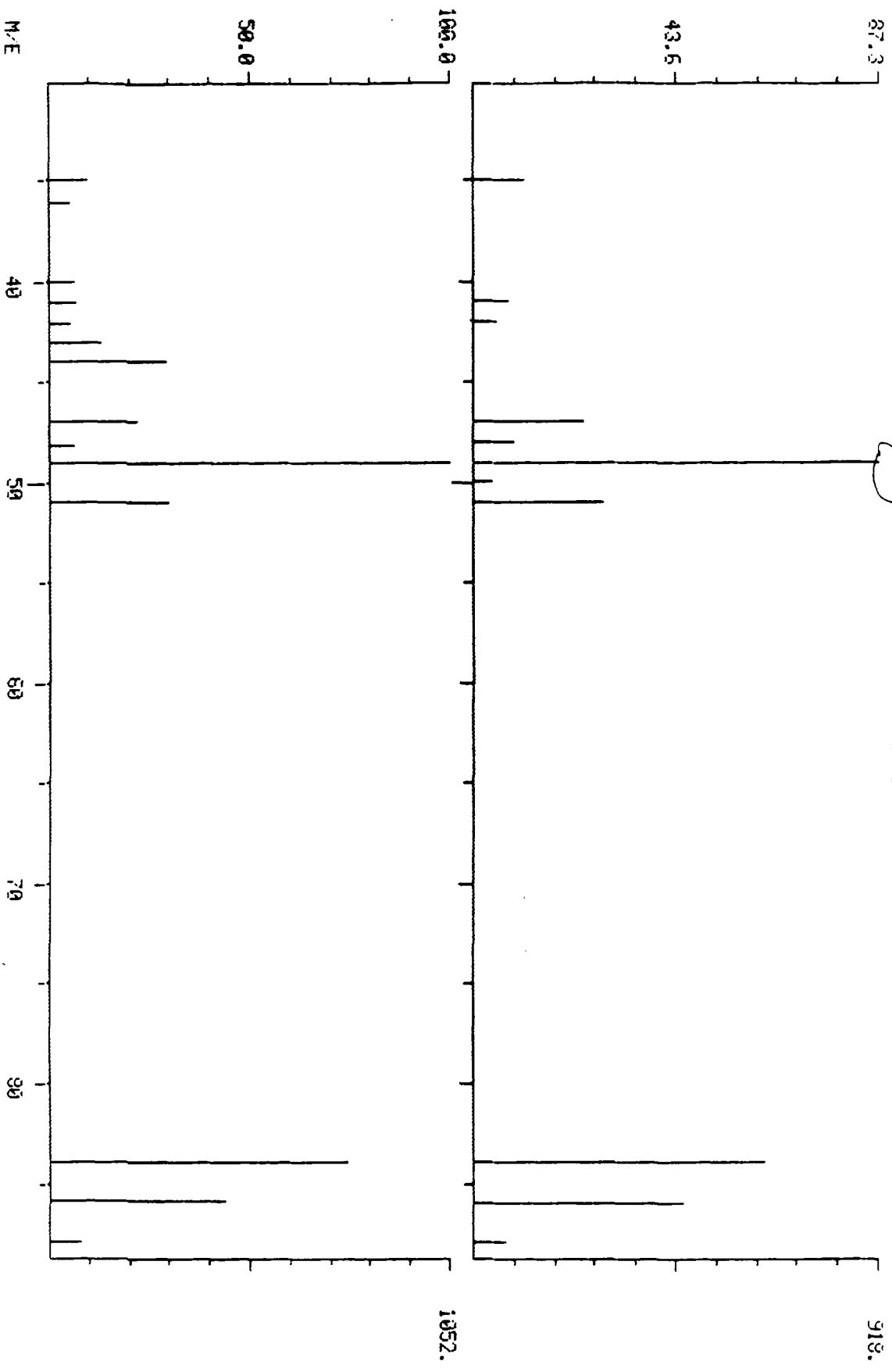
DUAL MASS SPECTRUM
12/29/83 20:10:00 + 2:58
SAMPLE: 100UL CC#309686 EPA#82018 CASE#18758 ON #13
ENHANCED (5 158 2N)

222

METHYLENE CHLORIDE (75-09-2) NA#16

COMPUchem LABS

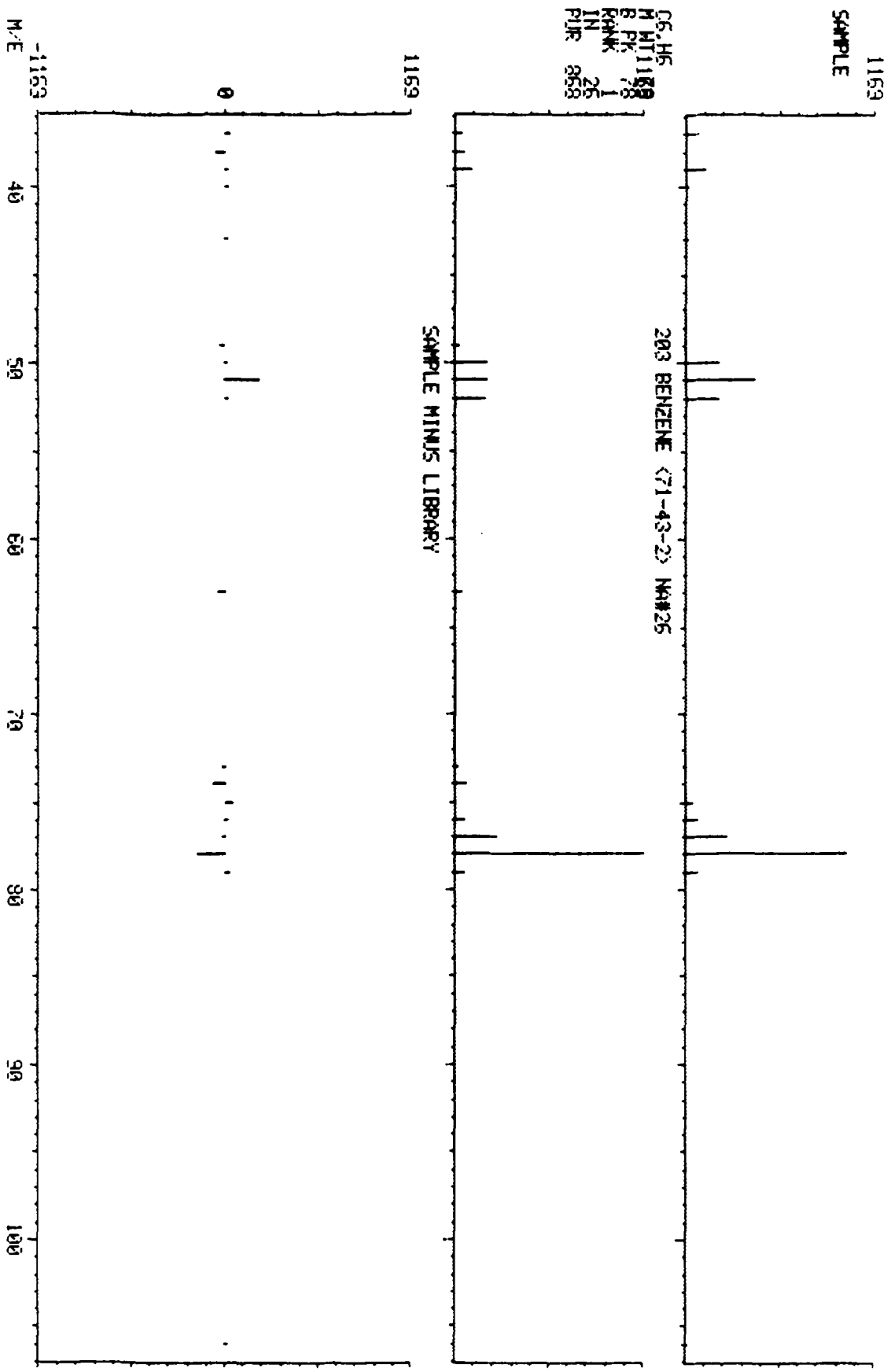
DATA: C5R09686B13 #238
BASE M/E: 49/ 49
RIC: 3043.7 3763.



COMPUCHEN LABS
LIBRARY SEARCH
12/29/89 20:10:00 + 5:55
SAMPLE: 100UL CC#309686 EPA#B2018 CASE#18755 ON #13
ENHANCED (5 158 2H 0T)

DATA: CSR09686813 # 474
BASE M/E: 78
RIC: 9439.

C6.H6
M.WT 118.0
B.Pk 78
KOKK
IN 26
PUR 8658



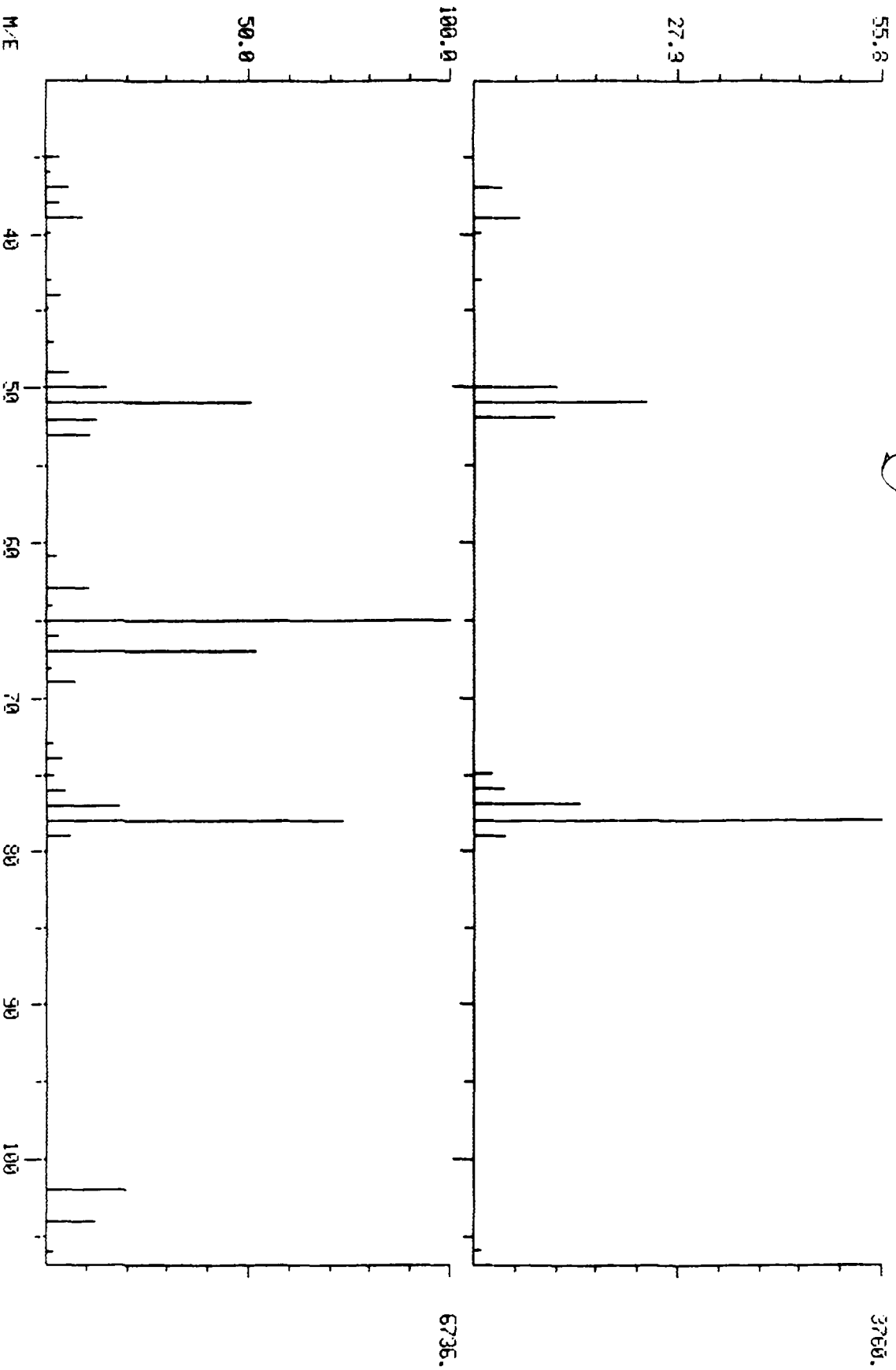
DUAL MASS SPECTRUM
12/23/89 20:10:00 + 5:55
SAMPLE: 100UL CC#309686 EPA#82018 CASE#18755 ON #13
ENHANCED (5 158 2N)

COMPUCHEM LABS
Ac of the Se Gb you like
203 BENZENE (71-43-2) NR#25

COMPUCHEM LABS

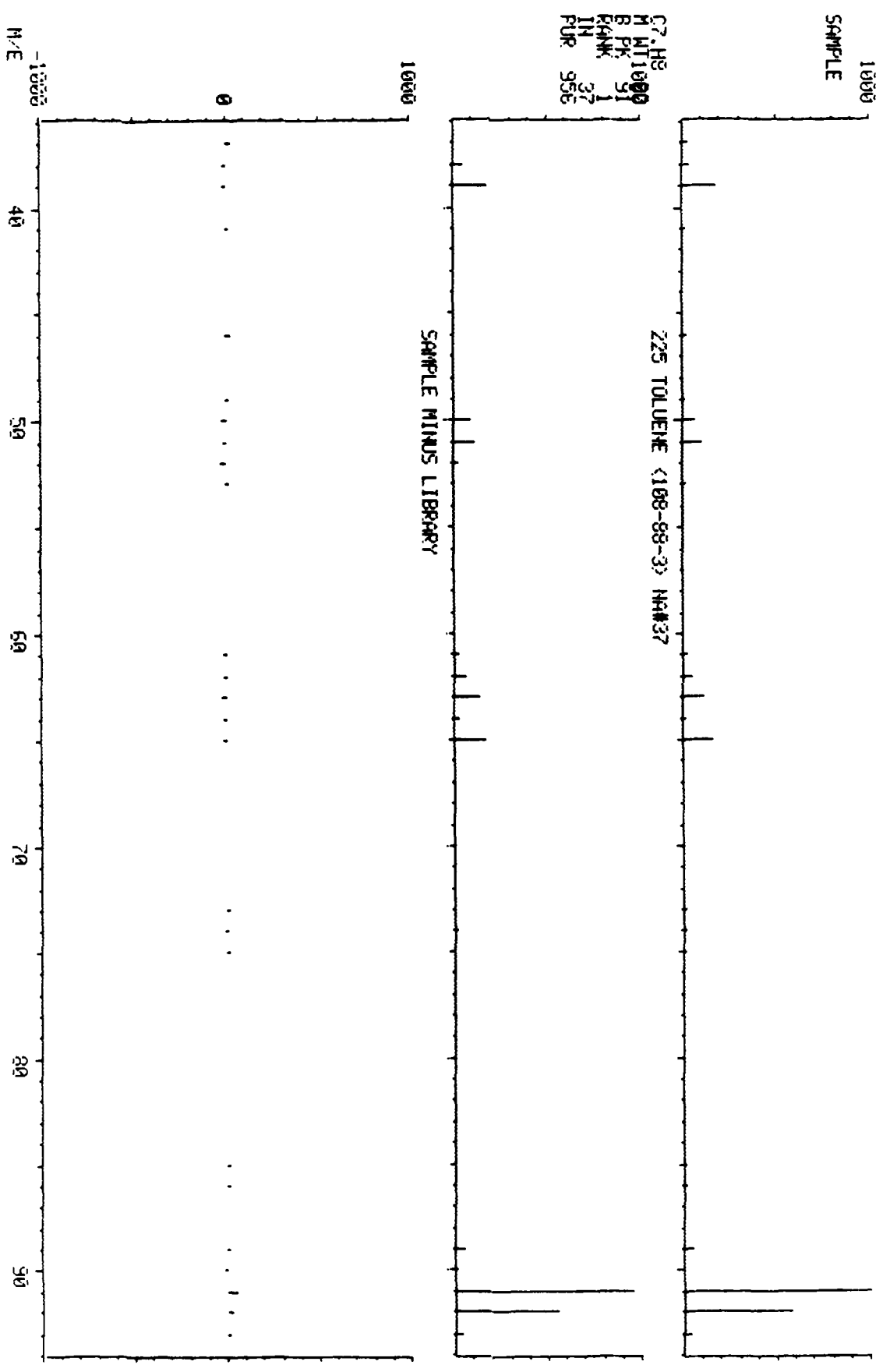
DATA: CSP09696813 #474

BASE M/E: 78/ 65
RIC: 9439. / 23887.



COMPUCHER LABS
 DATA: CSR09688613 # 723
 BASE M/E: 91
 LIBRARY SEARCH 12/29/89 20:10:00 + 9:02 R# 11296 G# 11296
 SAMPLE: 100UL CC#209688 EPANEZ01B CASE#18787 ON #13
 ENHANCED (5.158.2H.0T) RIC: 61759.

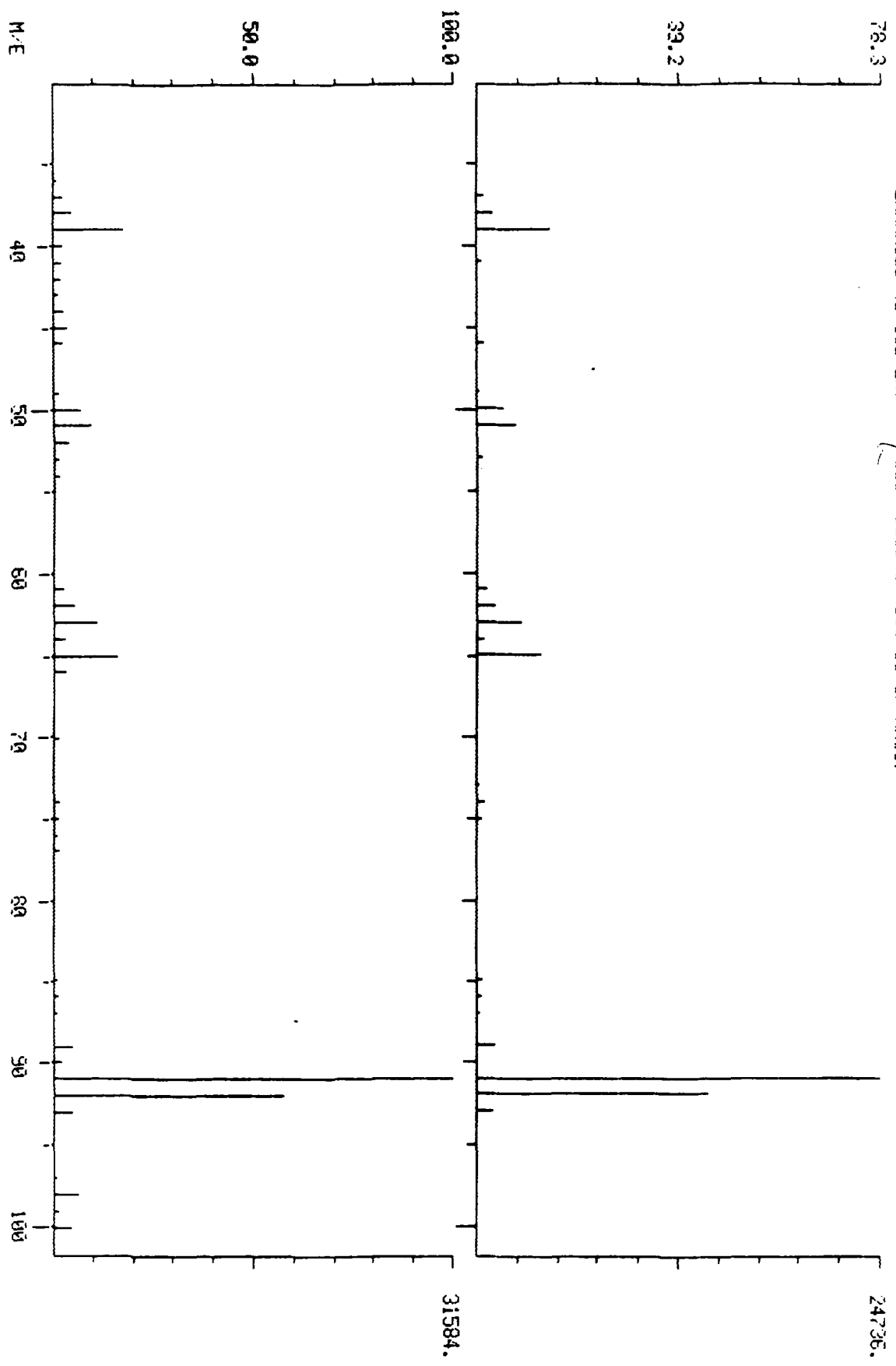
C7.H8
 H.M.I. 1000
 B.P.K. 91
 RANK 1
 IN 37
 PUR 996



DUAL MASS SPECTRUM
12/29/89 20:10:00 + 9:02
SAMPLE: 100UL CC#309686
ENHANCED (5 158 2N)

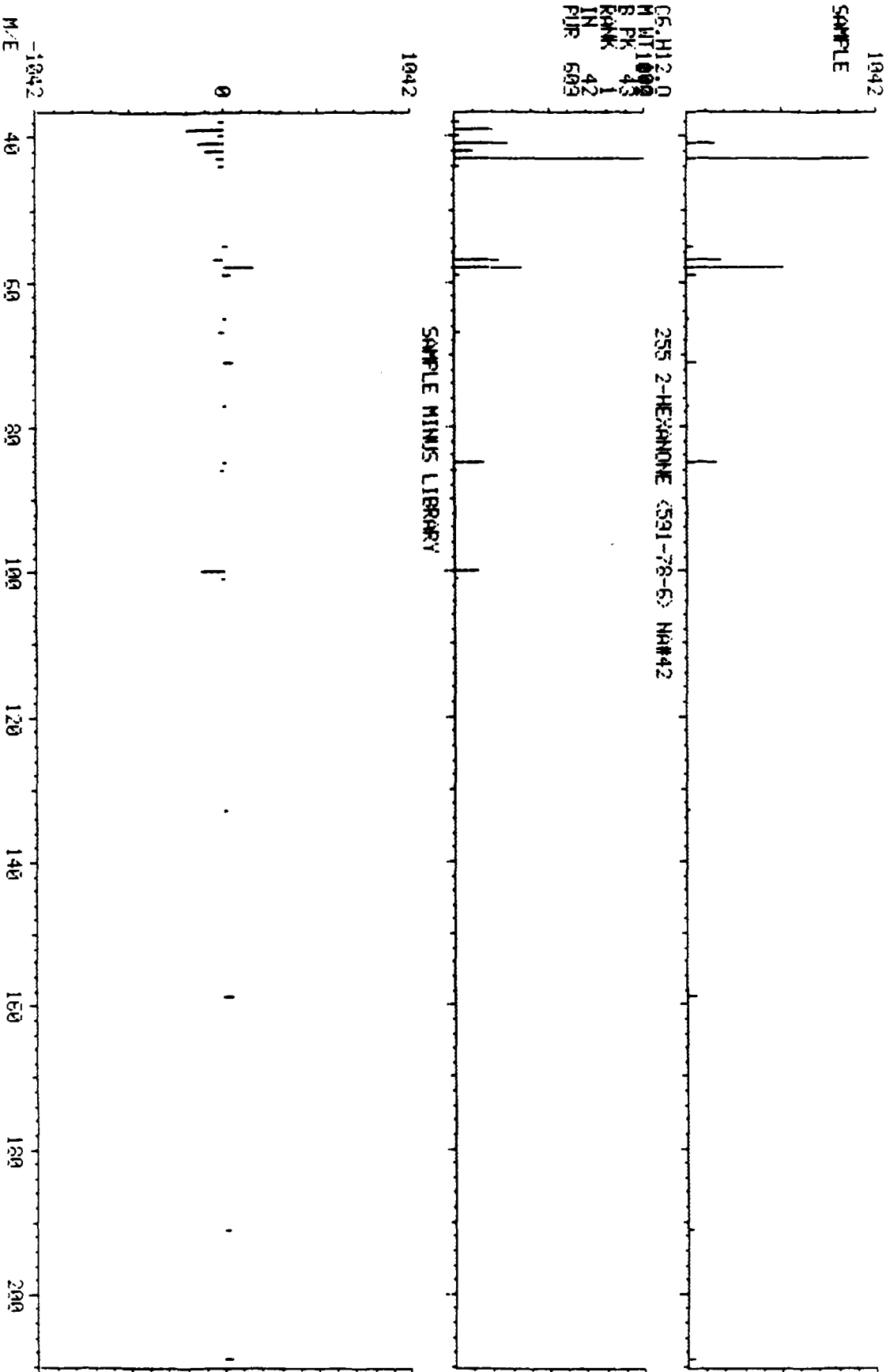
COMPUchem LABS
DATA: CSR09686813 #723
Revised by Sam Alder

BASE M/E: 91 / 91
RIC: 61951. / 91775.



COMPTON CHEM LABS
LIBRARY SEARCH
12/23/83 20:10:00 + 10:35
SAMPLE: 100UL CC#309686 EPA#B2018 CASE#18756 ON #13
ENHANCED (5 158 2N 0T)

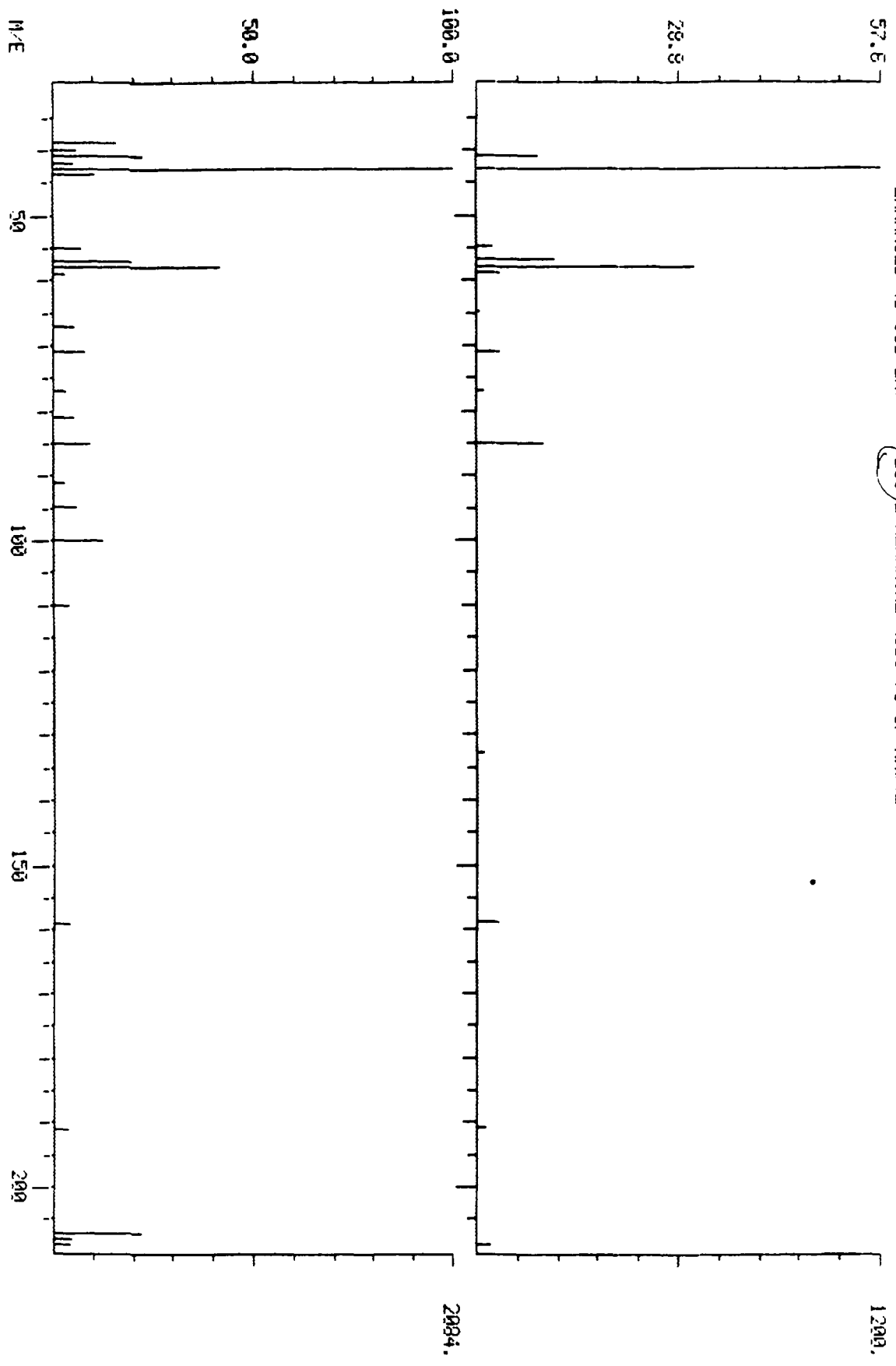
DATA: 05R03668B13 # 847
BASE M/E: 43
R10: 2815.



DUAL MASS SPECTRUM
12/29/89 20:10:00 + 10:35
SAMPLE: 100UL CC#309686
ENHANCED (5 158 2N)

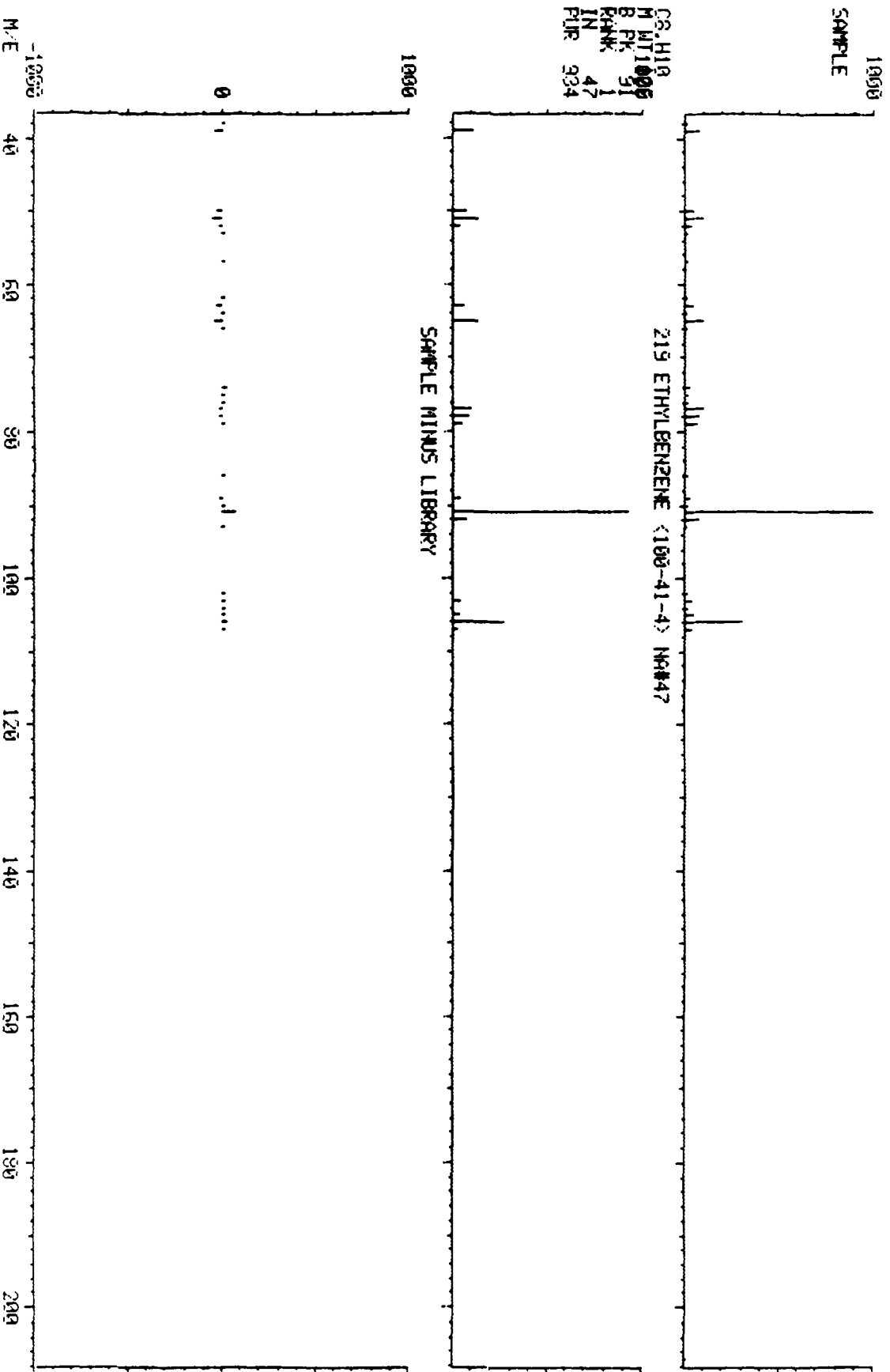
COMPUCHEM LABS
DATA: CSF09686B13 #847
255 2-HEXANONE (591-78-6) NA#42

BASE M/E: 43
RIC: 2815.7
6719.



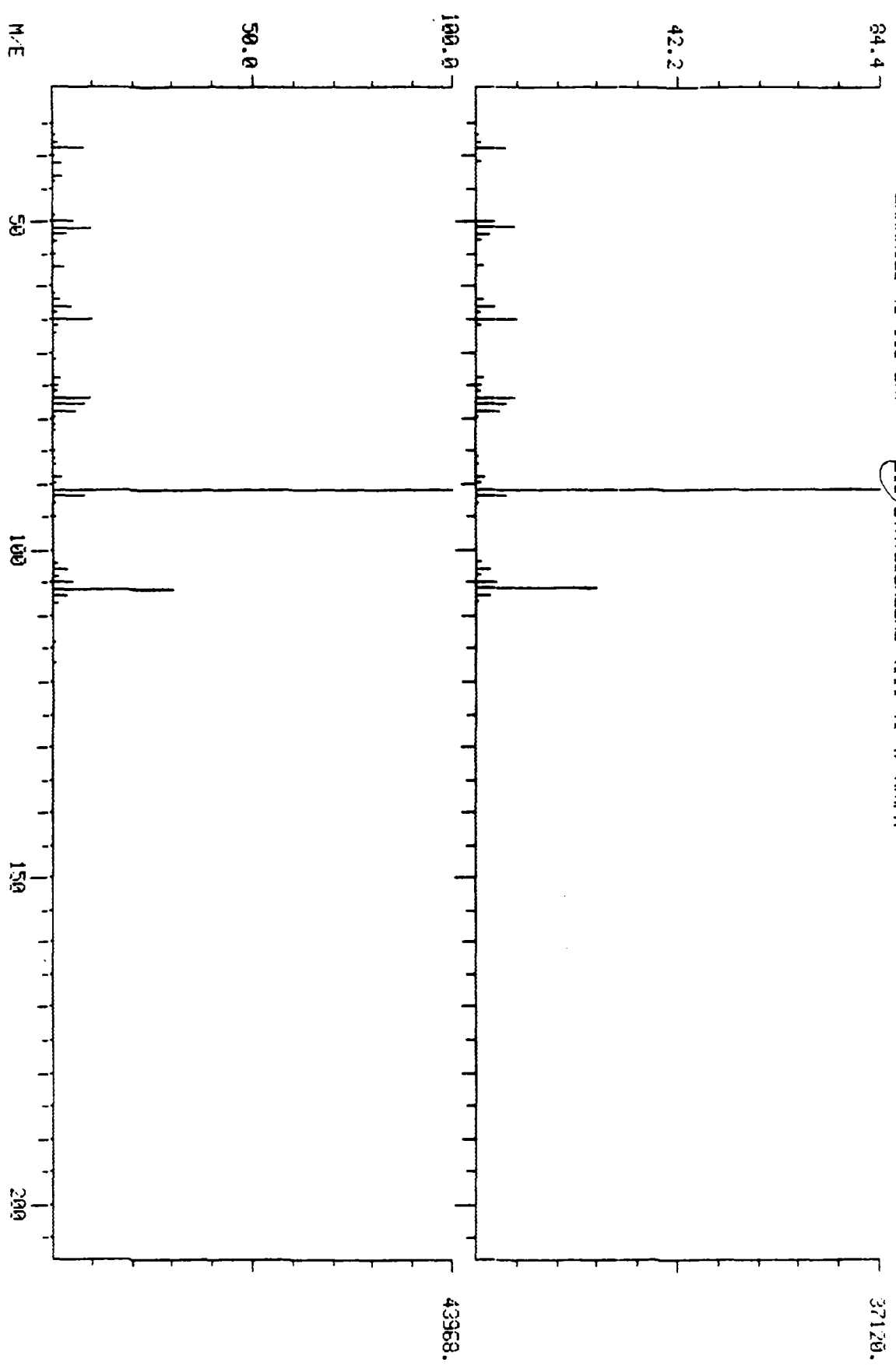
LIBRARY SEARCH
 12/29/89 20:10:00 + 11:46
 SAMPLE: 100UL C0309686 EPHABZ01B CASE#18763 ON #13
 ENHANCED (S 158 2N 0T)

COMPUchem LABS
 DATA: C0309686E13 # 942
 BASE M/E: 91
 RIC: 84863.



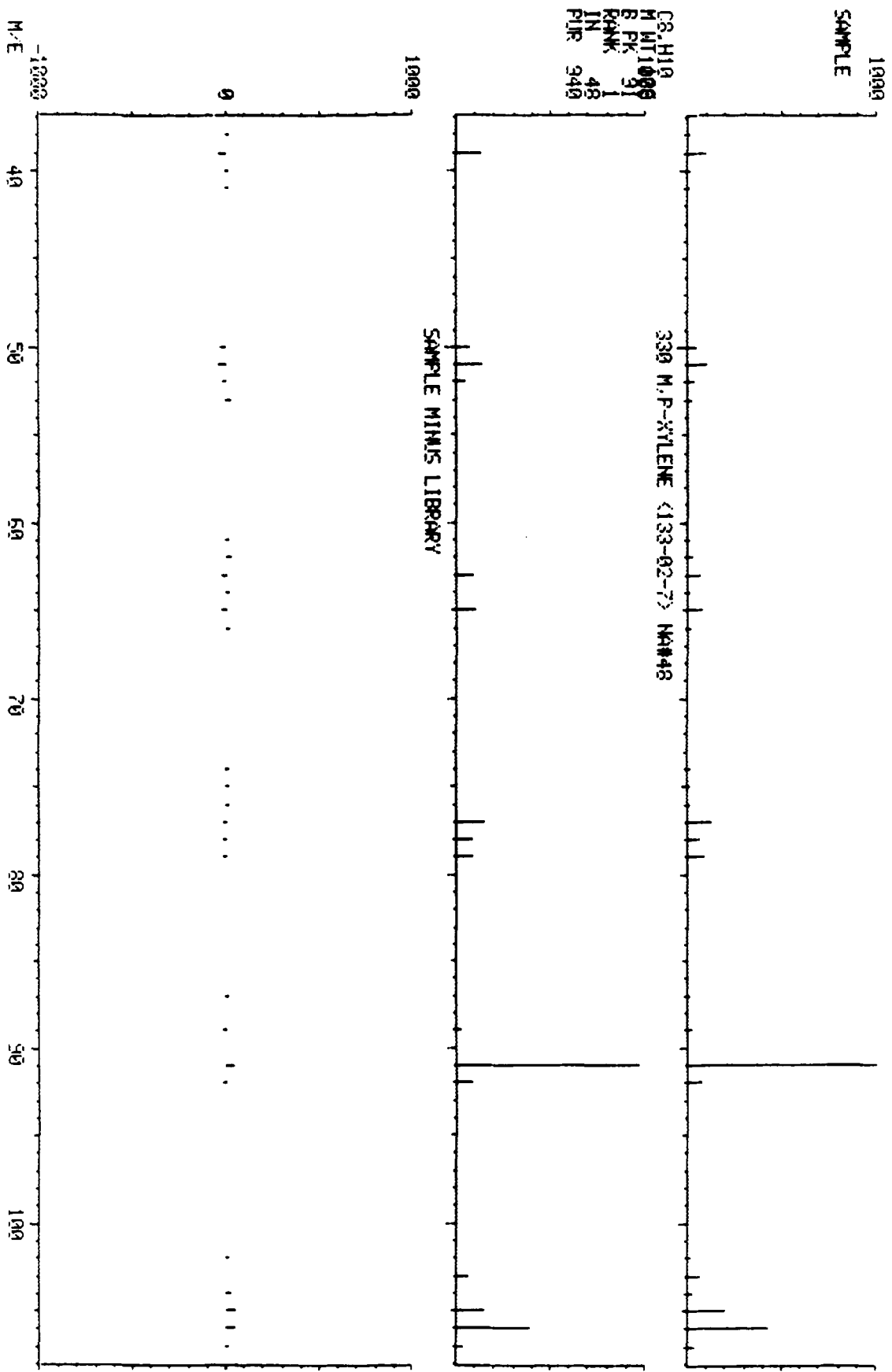
DUAL MASS SPECTRUM
12/29/89 20:10:00 + 11:46
SAMPLE: 100UL CC#309585 EPA#82018 CASE#18756 ON #13
ENHANCED (S 158 2N) (219)ETHYLBENZENE (100-41-4) NA#47

COMPUchem LABS
DATA: CSR09666813 #942
BASE M/E: 91 / 91
R/C: 86339. / 108927.



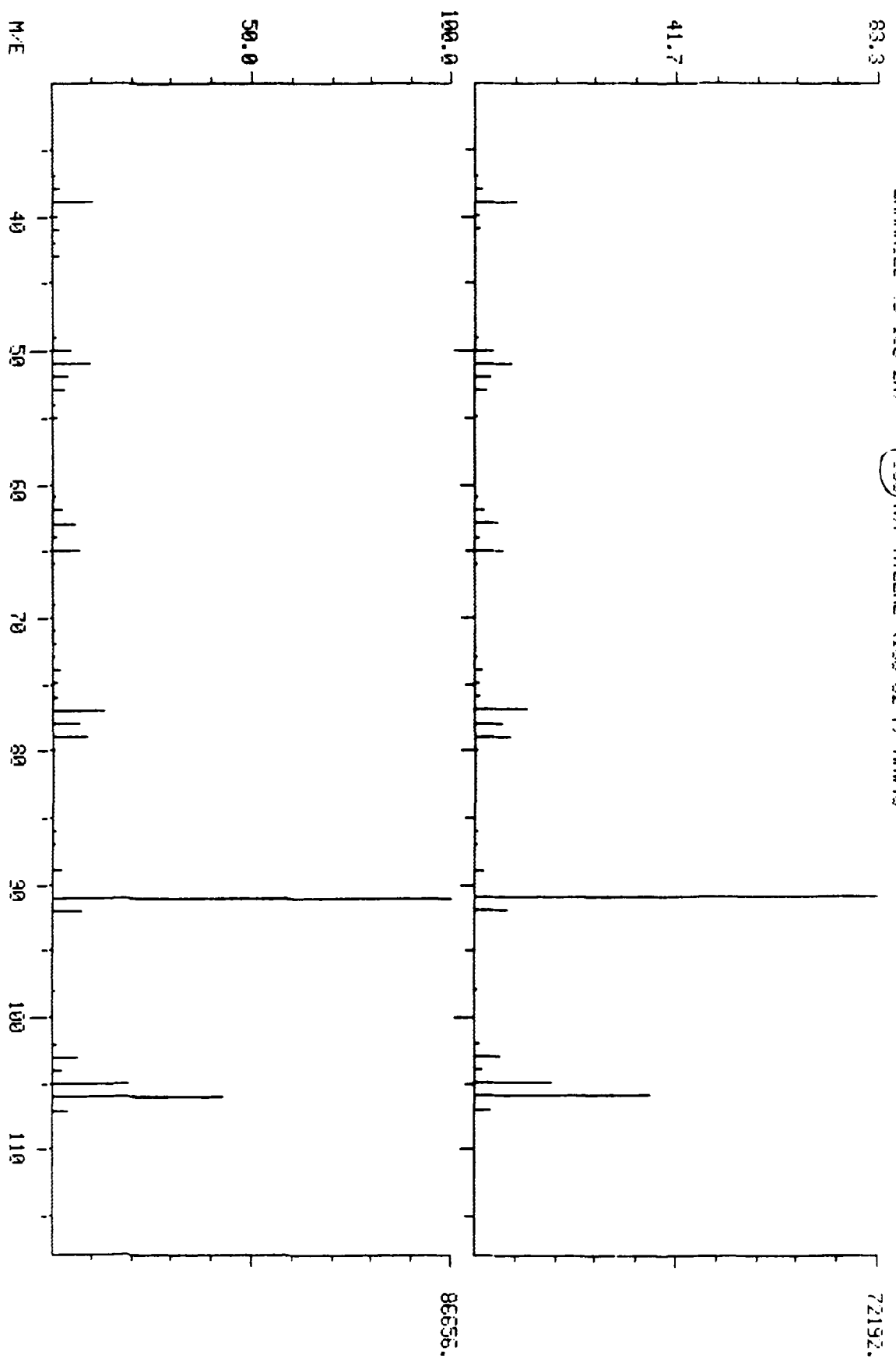
LIBRARY SEARCH
 12/29/89 20:10:00 + 12:00 RE
 SAMPLE: 100UL CCM309585 EPA#B2018 CASE#18755 ON #13
 ENHANCED (S 158 2H 0T)
 COMPUchem LABS
 DATA: CSR09658B13 # 360
 BASE M/E: 91
 RIC: 195071.

CG-HIQ
 M.WT 1000
 S.FK 31
 RANK 1
 IN 48
 PUR 940



DUAL MASS SPECTRUM
 12/29/89 20:10:00 + 12:00
 SAMPLE: 100UL CC#303685 EPA#B201B CASE#18756 ON #13
 ENHANCED (S 15R 2M) (330) M,P-XYLENE (133-02-7) NA#48

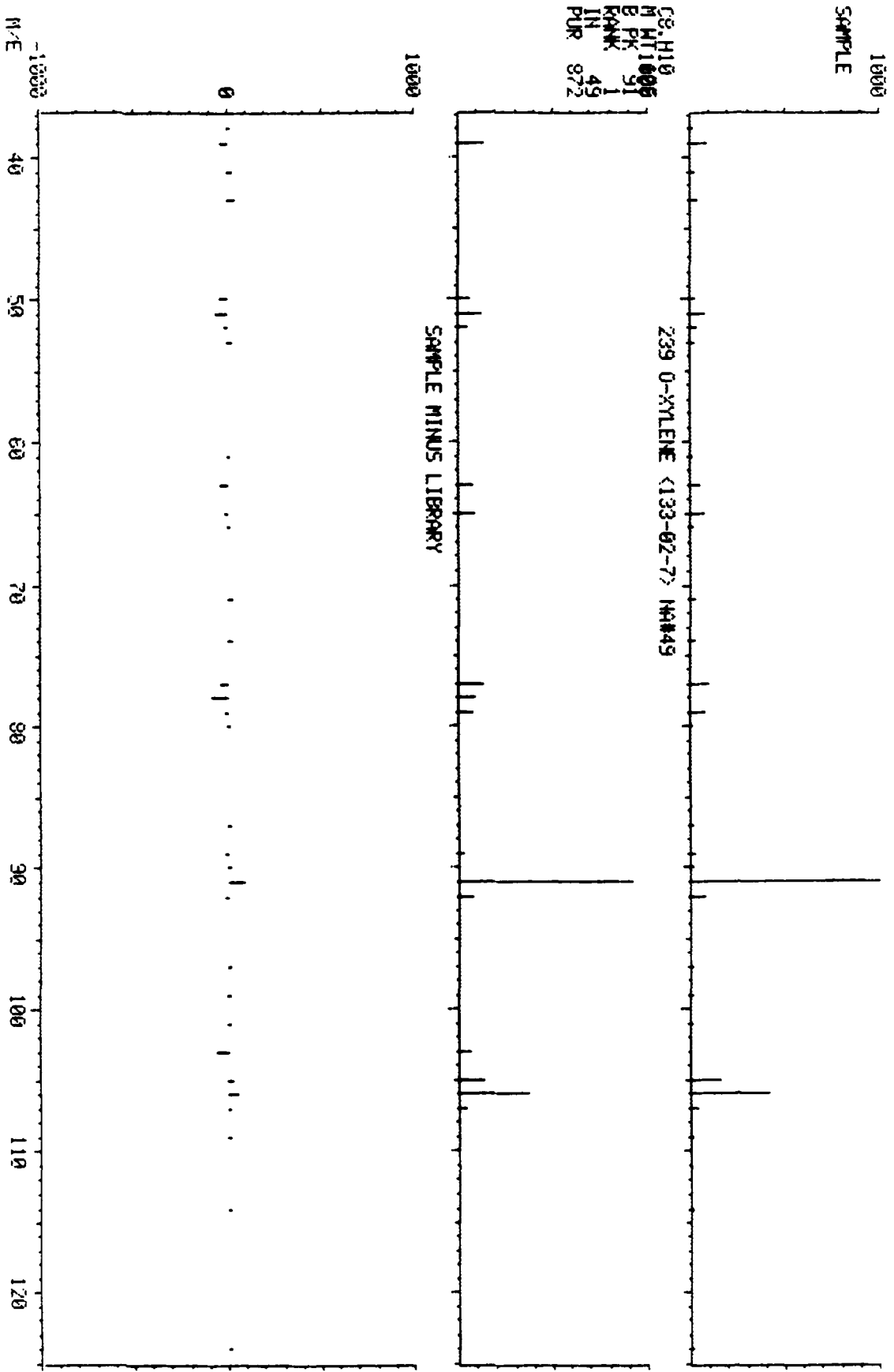
COMPUCHEM LABS
 DATA: CSR09696813 #960
 BASE M/E: 91/ 91
 RIC: 198399. / 243139.



LIBRARY SEARCH
 12/29/89 20:10:00 + 12:41
 SAMPLE: 100UL CC#309686 EPR#B201B CASE#18705 ON #13
 ENHANCED (S 158 2N 0T)

COMPUCHEN LABS
 DATA: CSR09686B13 #1015
 BASE M/E: 91
 RIC: 110079

CS: H10
 M: H1006
 B: PK 91
 RANK 49
 IN 1
 PUR 872

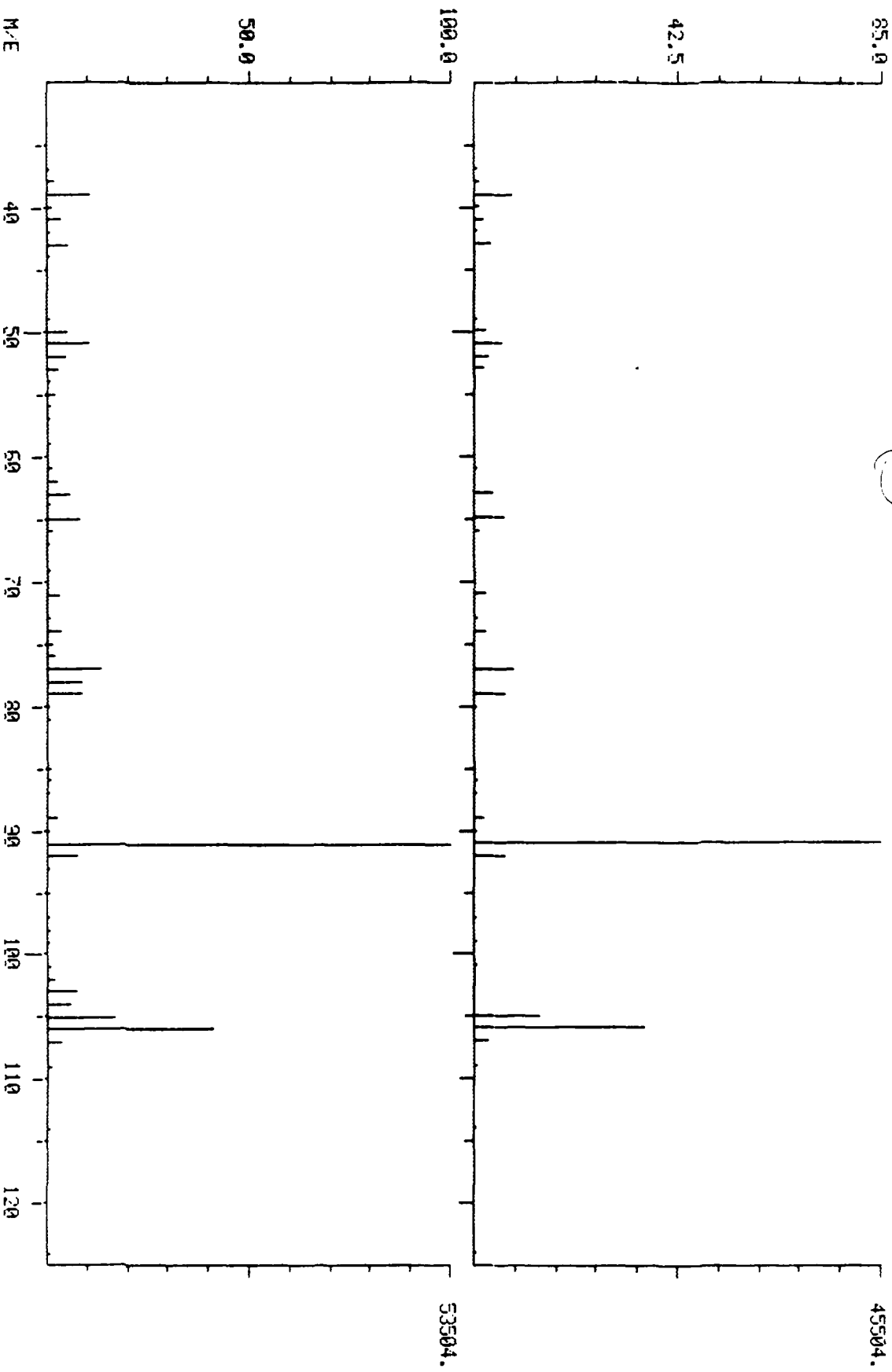


DUAL MASS SPECTRUM
12/29/89 20:10:00 + 12:41
SAMPLE: 100UL C0#309686 EPA#82018 CASE#18765 ON #13
ENHANCED (S 158 2N)

COMPUchem LABS
#6
CVS
SU
SAH
JALAB

DATA: C5R0666B13 #1015

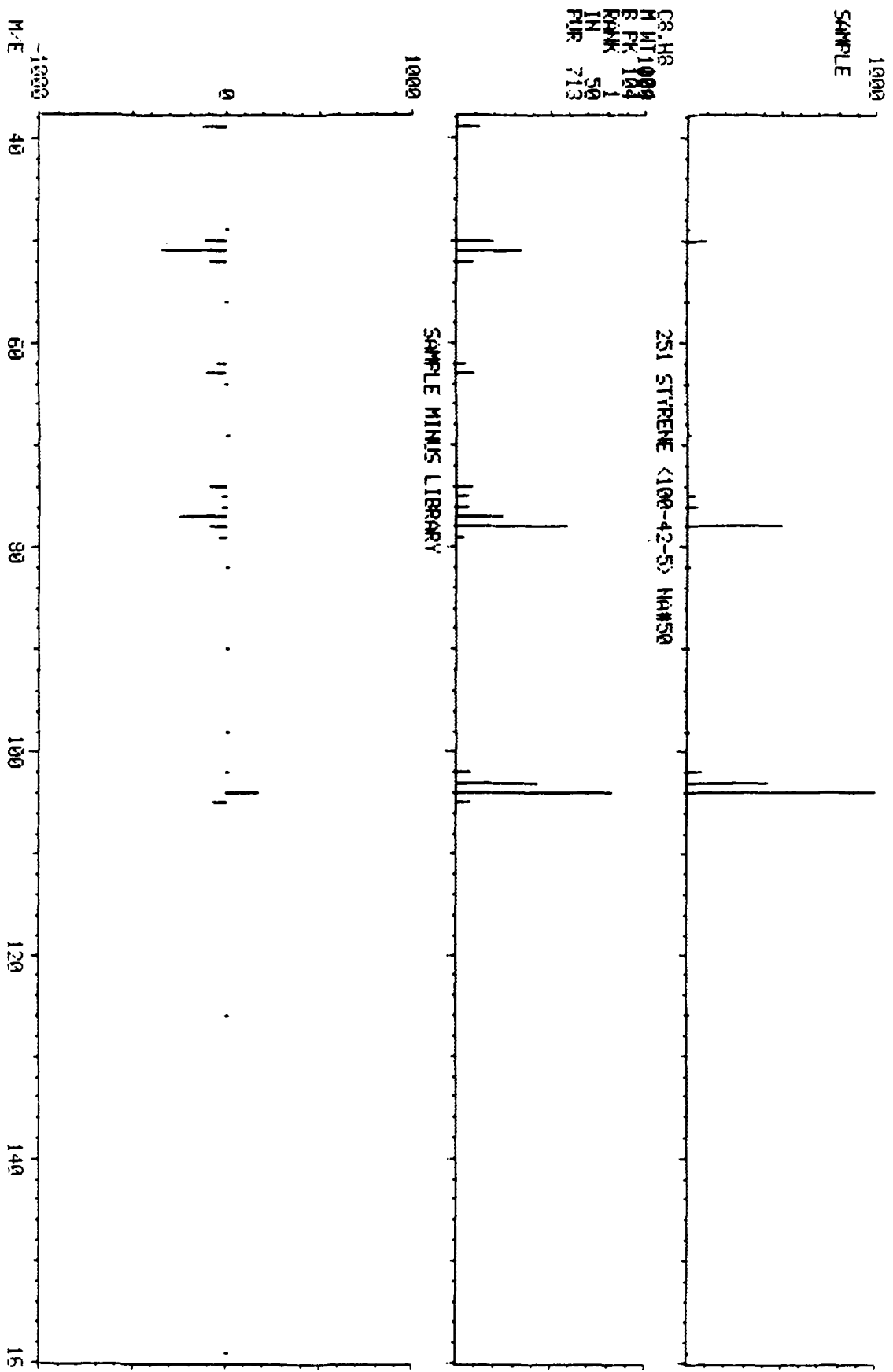
BASE M/E: 31/ 31
RIC: 112255. / 152815.



COMPUchem LABS
 LIBRARY SEARCH
 12/29/89 20:10:00 + 12:45
 SAMPLE: 100UL CC#309686 EPA#B201B CASE#18785 DN #13
 ENHANCED (5 158 2N 0T)

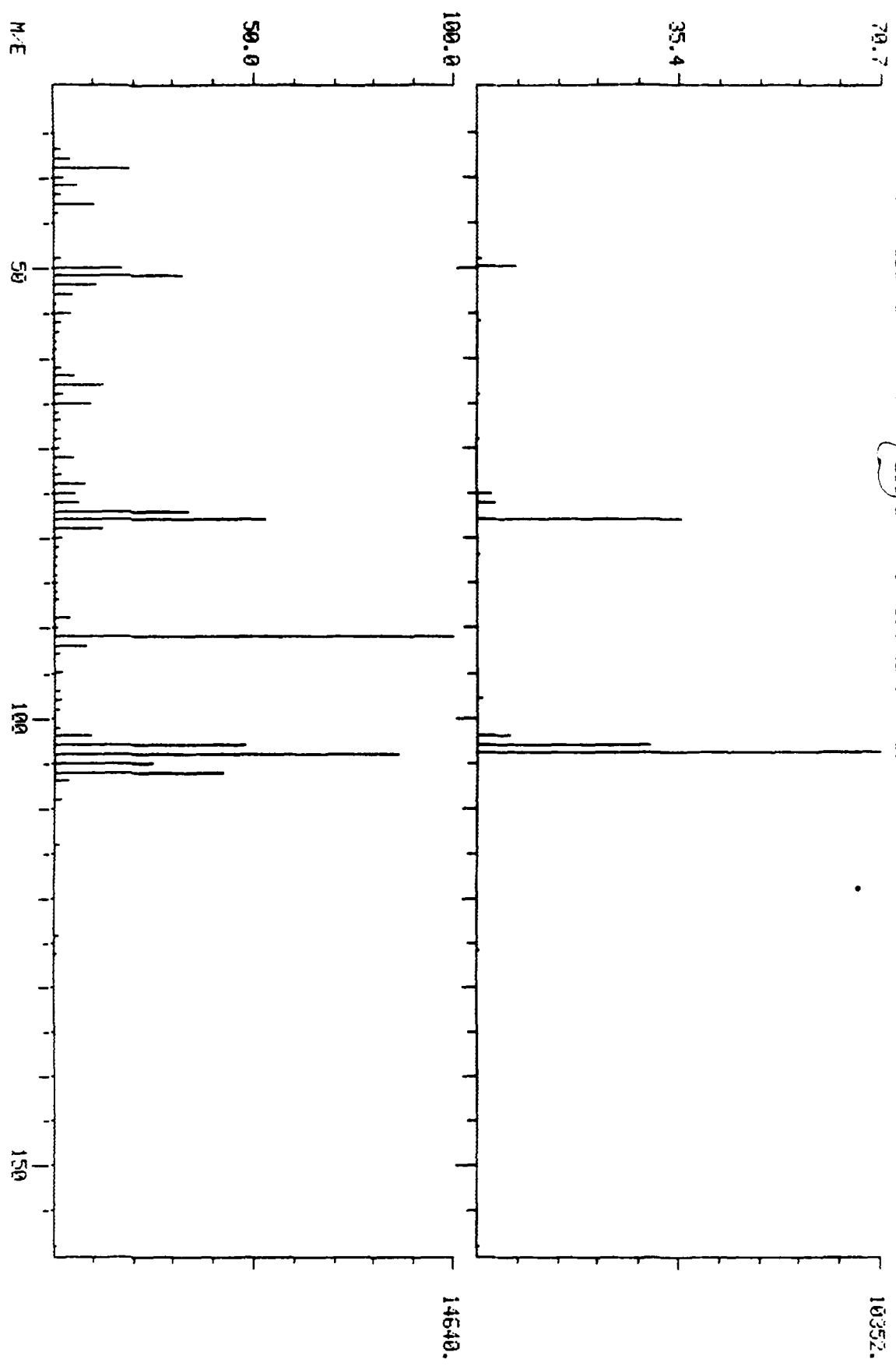
DATA: C5R09686B13 #1021
 BASE M/E: 104
 RIC: 23231.

C8.H8
 M WT 1004
 R PK 104
 RANK 1
 IN 50
 PUR 713



DUAL MASS SPECTRUM
12/29/89 20:10:00 + 12:45
SAMPLE: 100UL CC#309686 EPA#82018 CASE#18755 ON #13
ENHANCED (S 158 2N) 251 STYRENE (100-42-5) NAME59

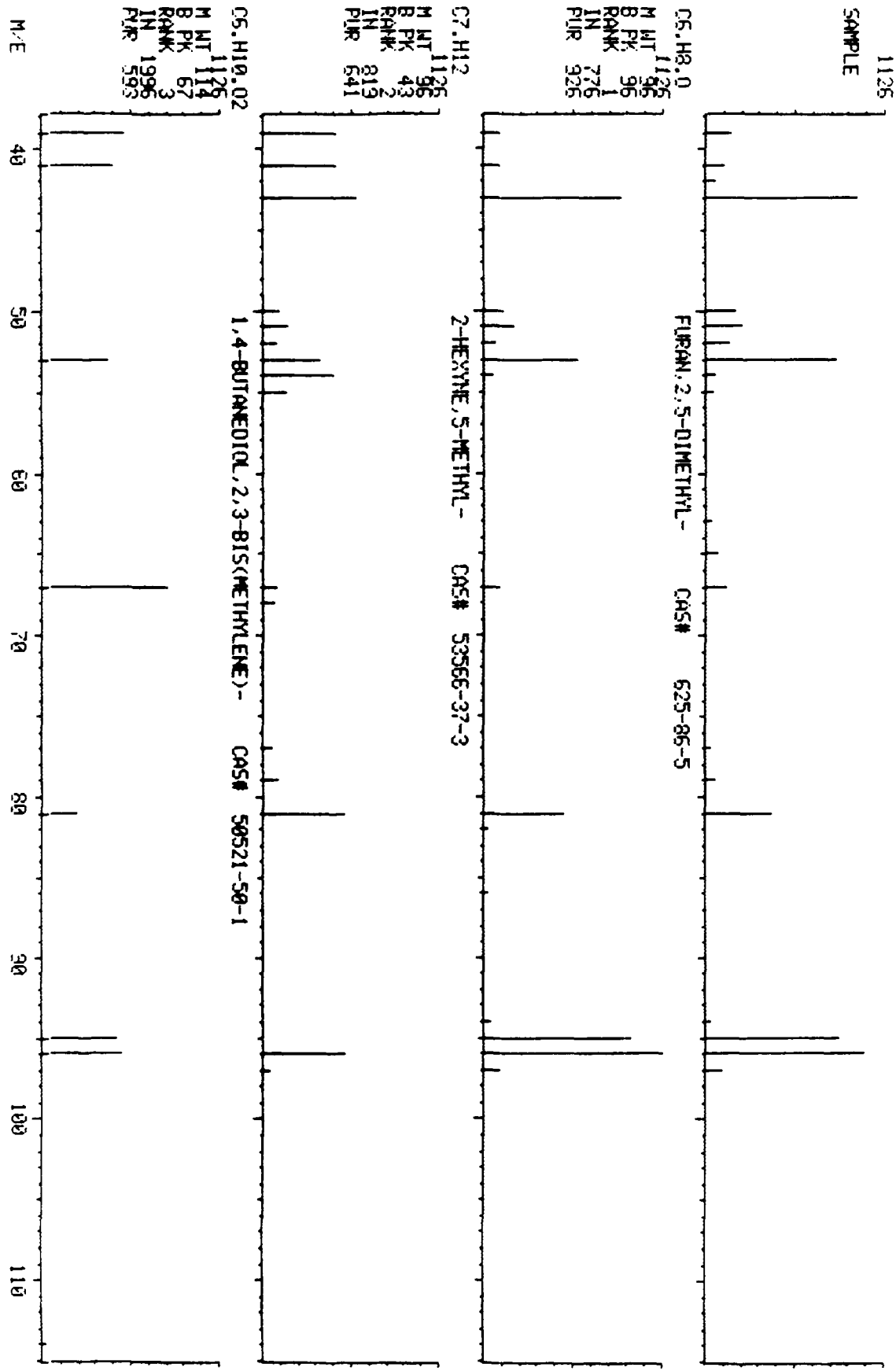
COMPUchem LABS
DATA: CSP09686813 #1021 BASE M/E: 104 / 91
RIC: 23231. / 91519.



LIBRARY SEARCH
 12/29/89 20:10:00 + 7:17
 SAMPLE: 100UL CC#309585 EPA#B201B CASE#18768 ON #13
 ENHANCED (S 158 2N 0T)

COMPUchem LABS
 DATA: CSR09686B13 # 583

BASE M/E: 96
 RIC: 13247.



LIBRARY SEARCH
 12/29/89 20:10:00 + 10:01
 SAMPLE: 100UL CC#309686 EPA#B201B CASE#18705 ON #13
 ENHANCED (S 158 2N 0T)

COMPUCHEN LABS
 DATA: CSR09686B13 # 802
 BASE M/E: 95
 RIC: 23615.

1000
 SAMPLE

05.H5.02
 M UT 1000
 B PK 95
 RANK 1
 IN 1598
 PUR 811

ETHANONE,1-(2-FURANYL)- CAS# 1192-62-7

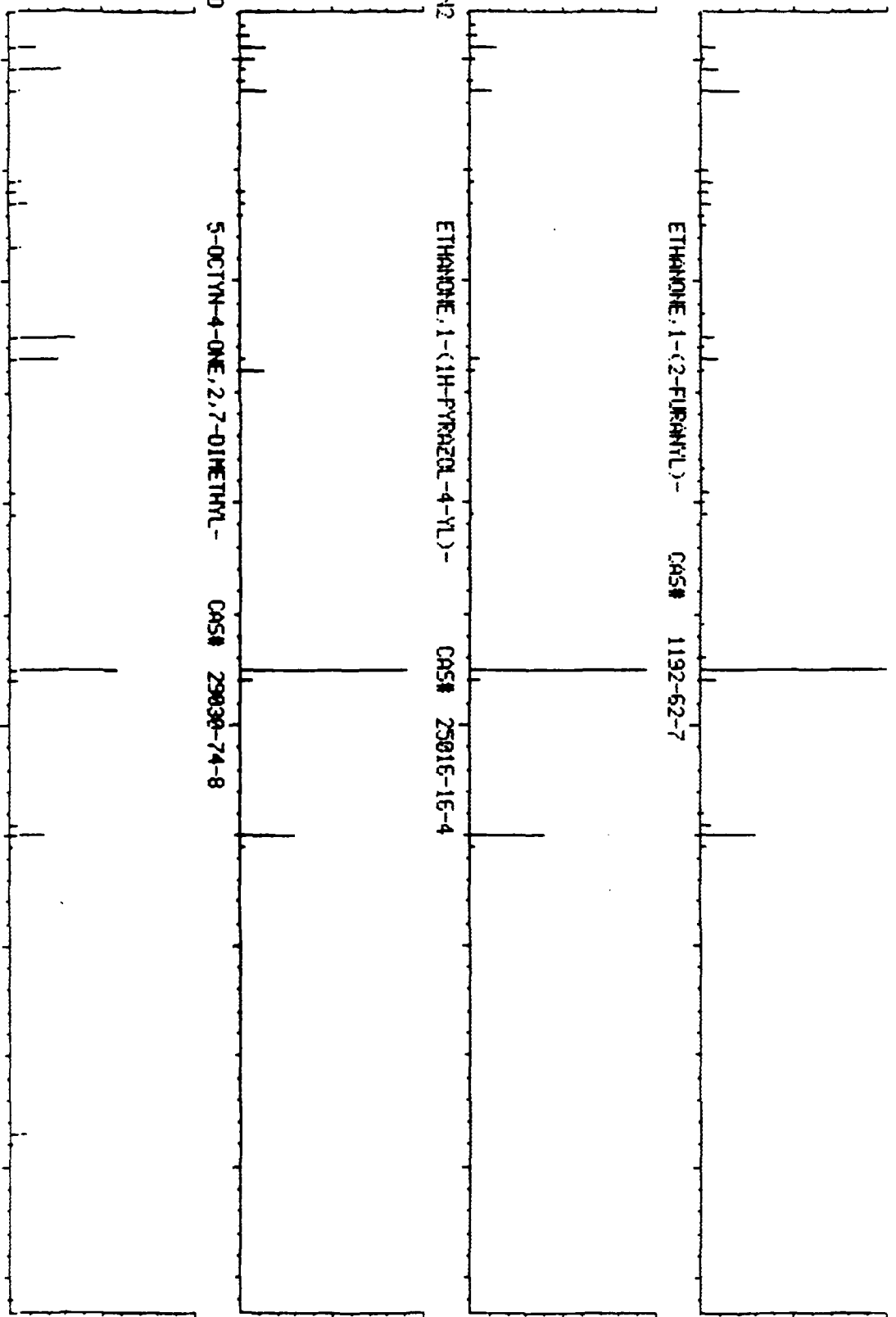
05.H5.0.N2
 M UT 1000
 B PK 95
 RANK 2
 IN 1588
 PUR 778

ETHANONE,1-(1H-PYRAZOL-4-YL)- CAS# 25016-16-4

C10.H16.0
 M UT 1000
 B PK 95
 RANK 3
 IN 6245
 PUR 735

5-OCTYN-4-ONE,2,7-DIMETHYL- CAS# 29030-74-8

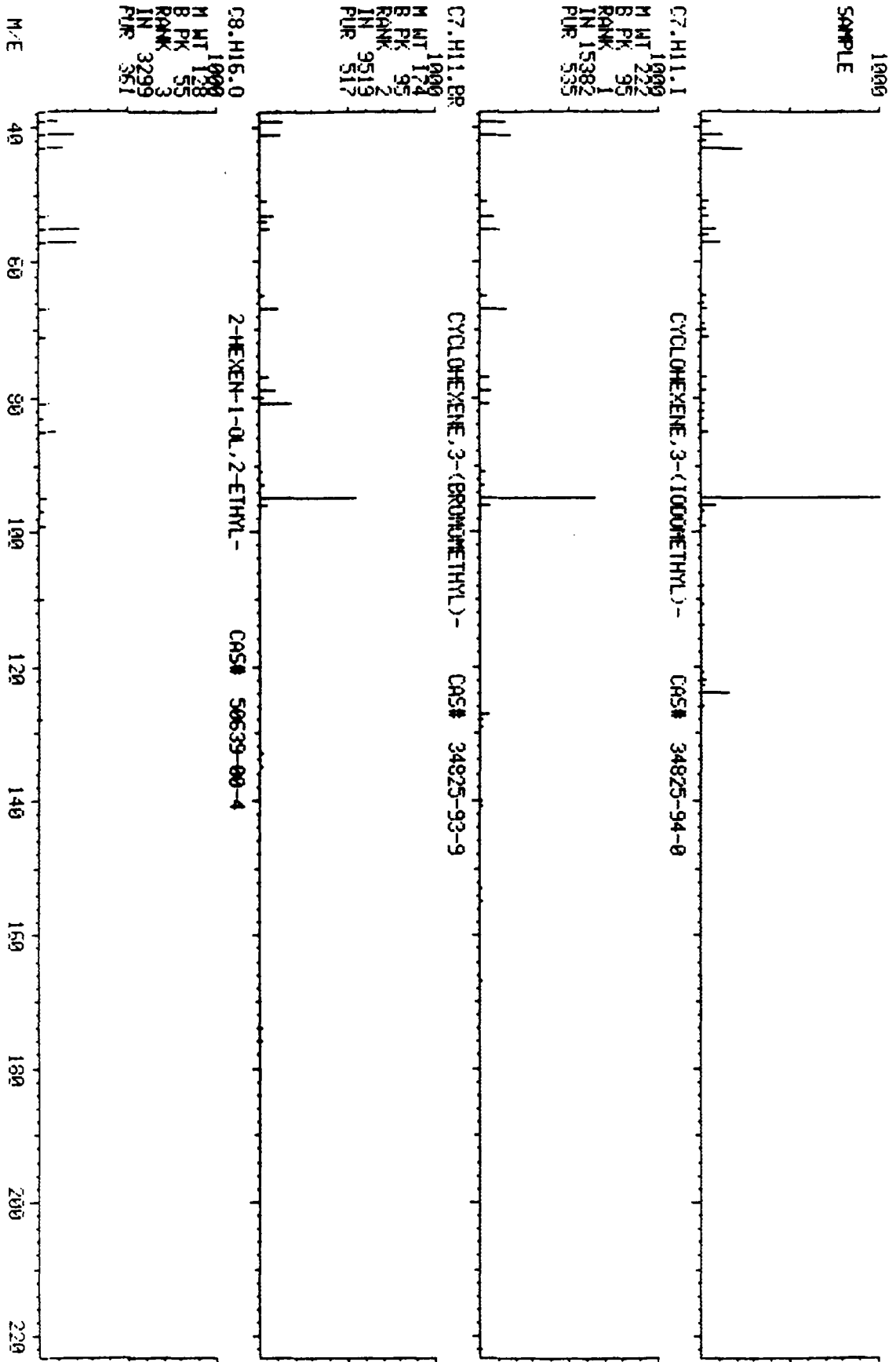
M/E 40 60 80 100 120 140



LIBRARY SEARCH
 12/29/89 20:10:00 + 12:20
 SAMPLE: 100UL CC#309686 EPA#B2018 CASE#18756 ON #13
 ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

DATA: CSR09686B13 # 987
 BASE M/E: 95
 RIC: 25631.

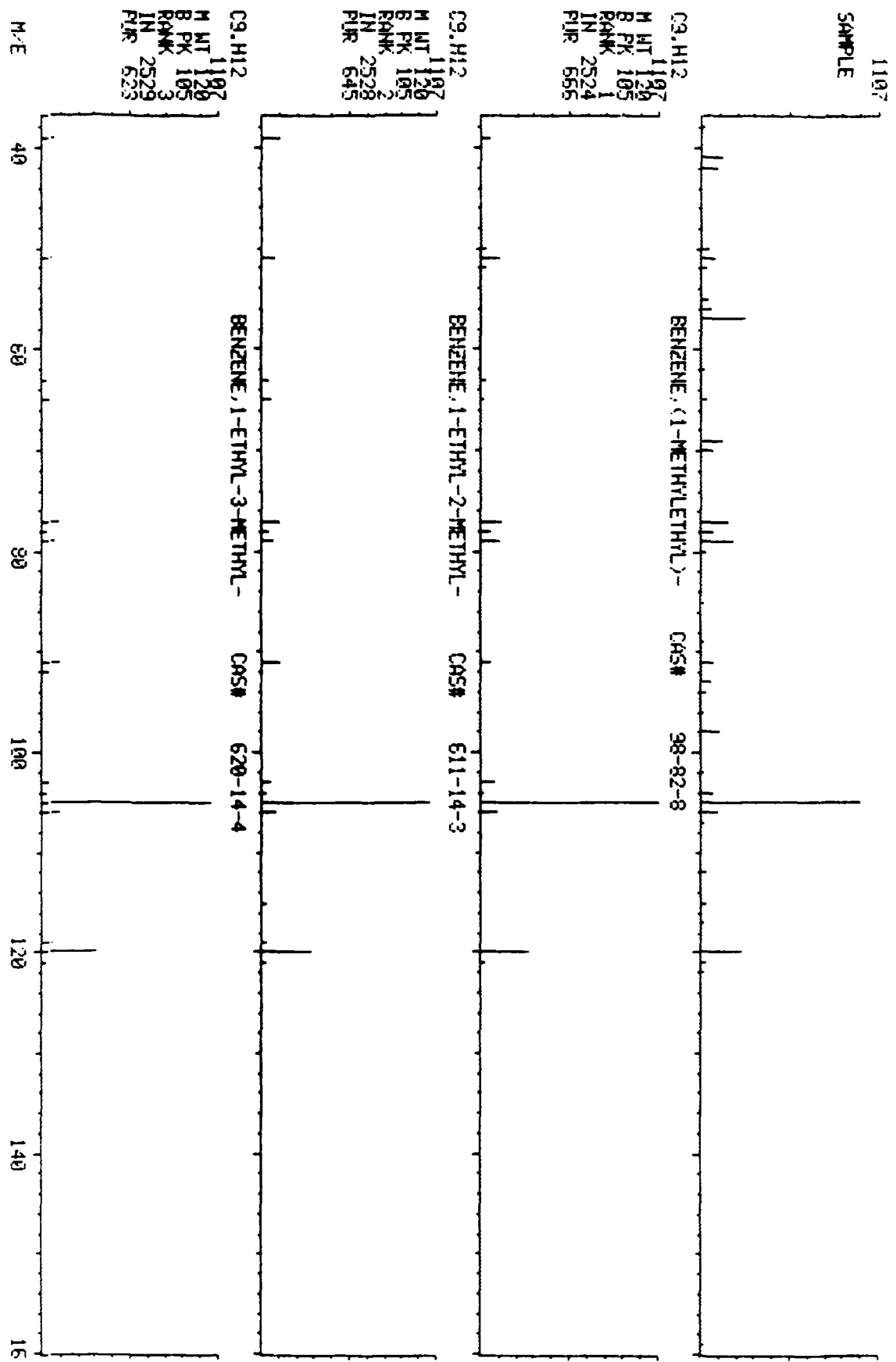


LIBRARY SEARCH
 12/29/89 20:10:00 + 13:26
 SAMPLE: 100UL CC#393685 EPA#B2018 CASE#18765 ON #13
 ENHANCED (S 158 2N 0T)

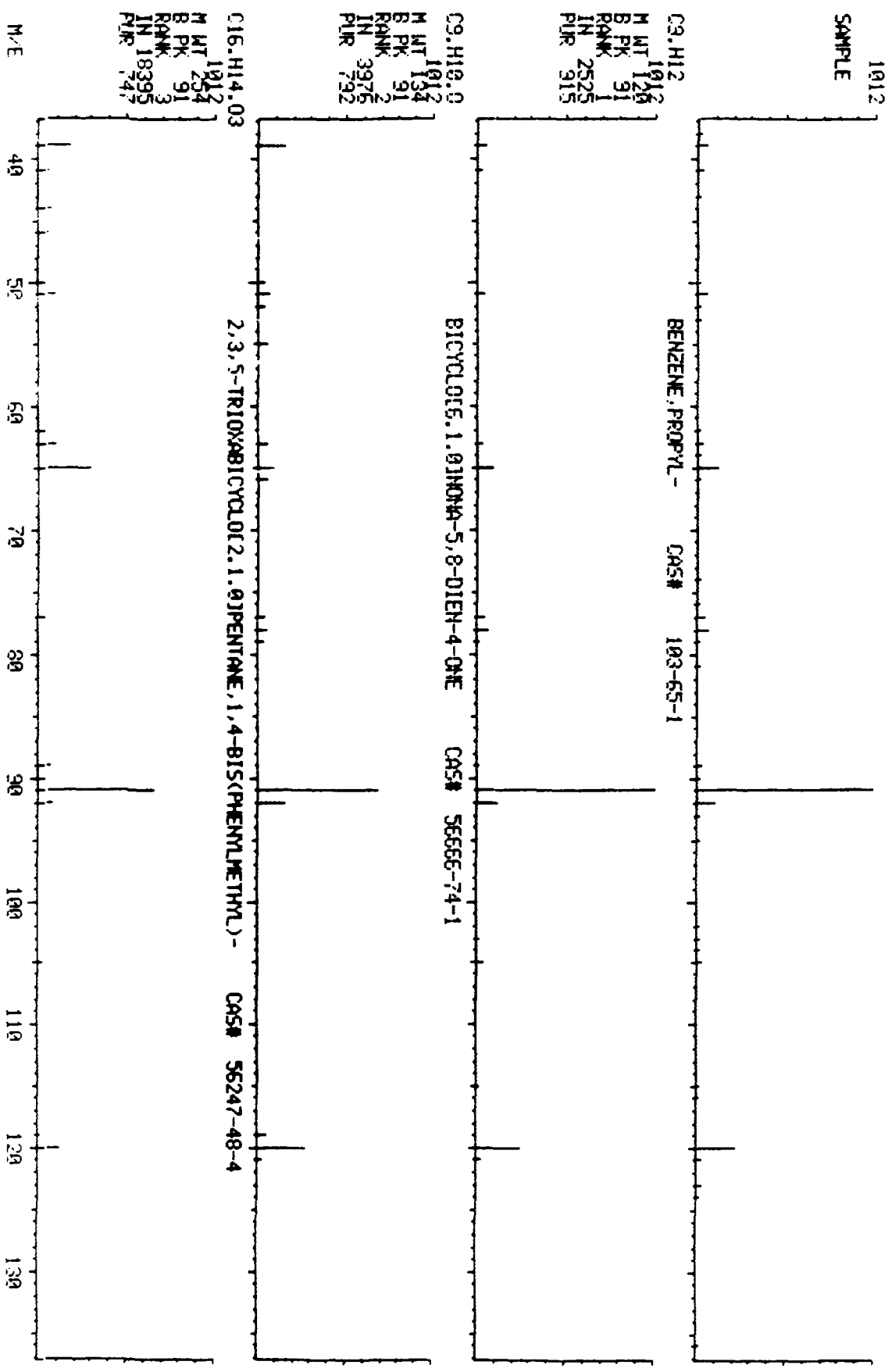
COMPUCHEM LABS

DATA: C5R03686613 #1075

BASE M/E: 105
 RIC: 13935.



COMPUchem LABS
 LIBRARY SEARCH
 12/29/89 20:10:00 + 14:13
 SAMPLE: 100UL CC#309685 EPA#82018 CASE#18756 ON #13
 ENHANCED (5 158 2N 9T)
 DATA: CSR09686813 #1138
 BASE M/E: 91
 RIC: 47295.



LIBRARY SEARCH
12/29/89 20:10:00 + 14:27
SAMPLE: 100UL CC#309586 EPA#B201B CASE#18756 ON #13
ENHANCED (5 158 2N 0T)

COMPUCHER LABS

DATA: C5R09586B13 #1156

BASE M/E: 105
R1C: 150527.

1040
SAMPLE

C9.H12
M NT 1040
B PK 105
RANK 1
IN 2524
PUR 896

BENZENE, (1-METHYLETHYL)- CAS# 98-82-8

C9.H12

M NT 1040
B PK 105
RANK 2
IN 2528
PUR 896

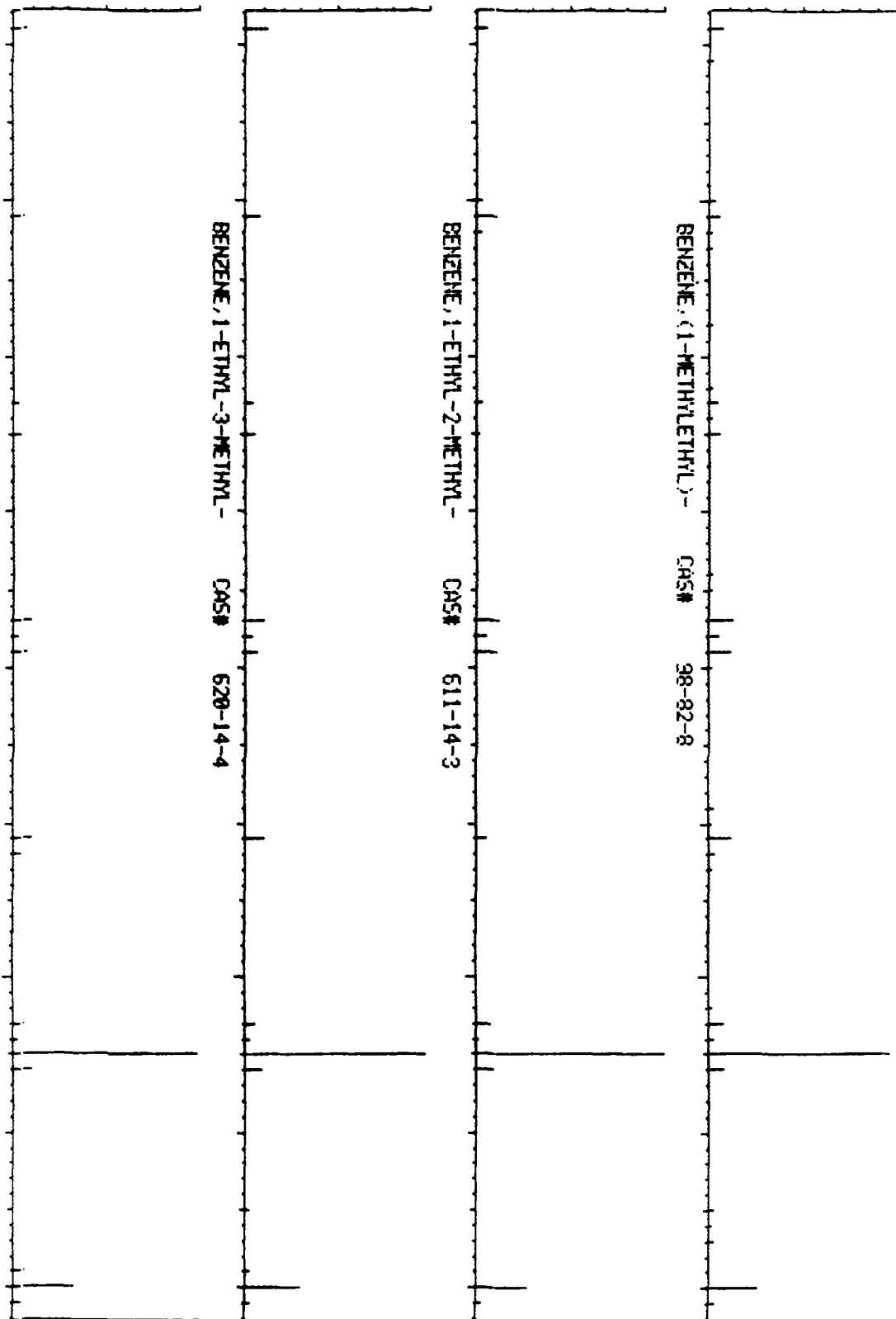
BENZENE, 1-ETHYL-2-METHYL- CAS# 611-14-3

C9.H12

M NT 1040
B PK 105
RANK 3
IN 2529
PUR 879

BENZENE, 1-ETHYL-3-METHYL- CAS# 620-14-4

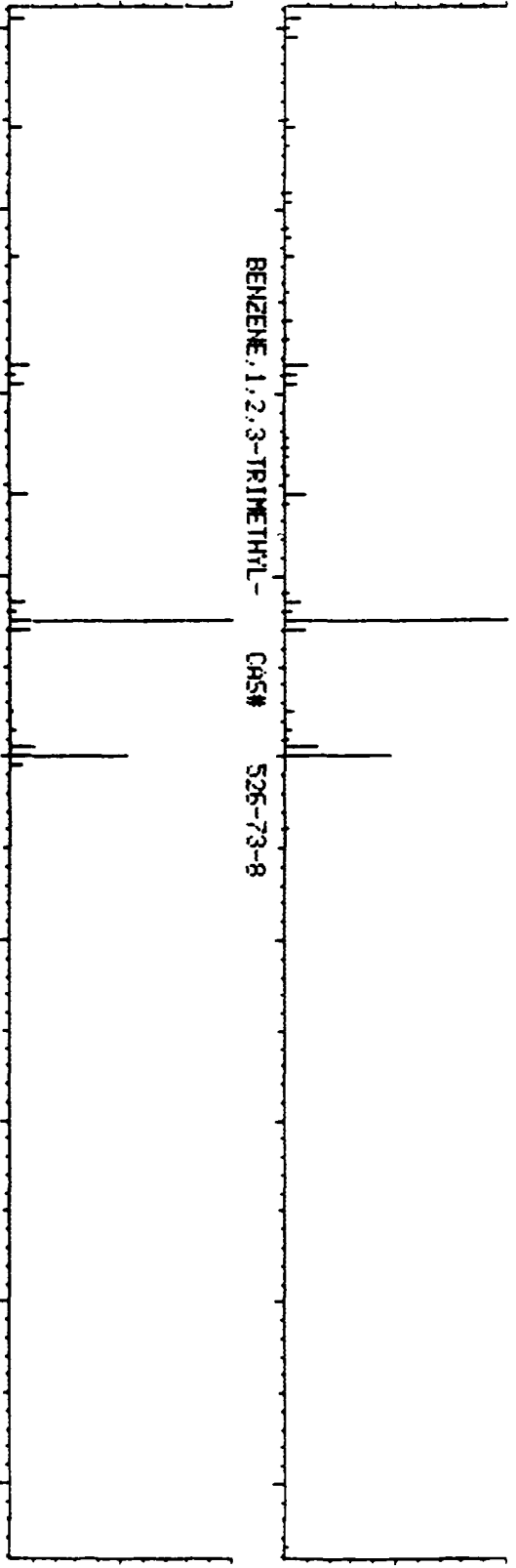
M/E 40 50 60 70 80 90 100 110 120



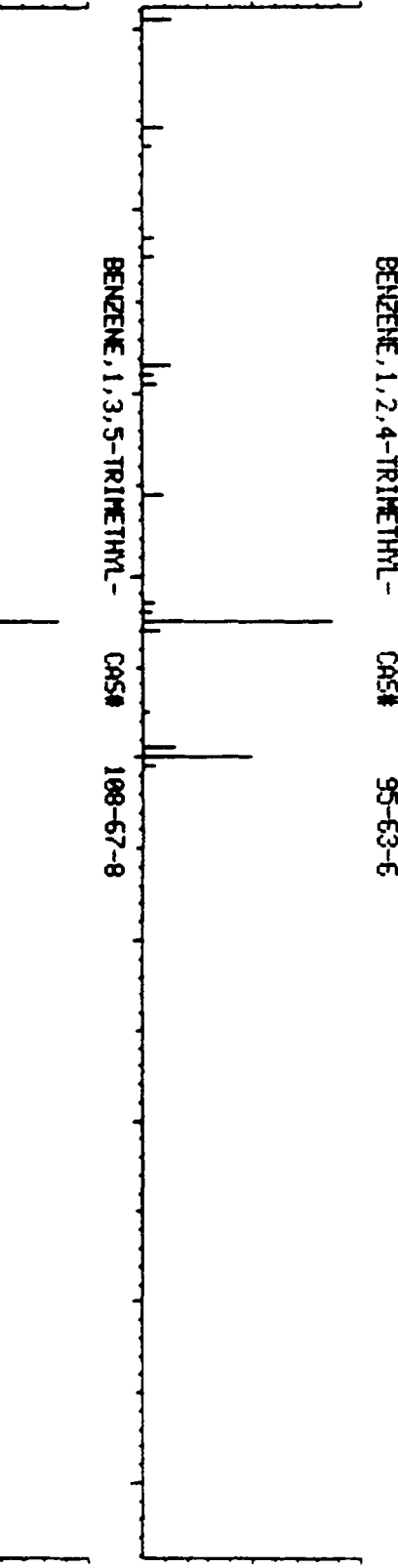
COMPUCHEN LABS
 LIBRARY SEARCH
 12/29/89 20:10:00 + 14:37
 SAMPLE: 100UL CC#399686 EPA#B201B CASE#18756 ON #13
 ENHANCED (5 158 2N 0T)
 DATA: C5R09686B13 #1169
 BASE M/E: 105
 RIC: 39615.

1002
 SAMPLE

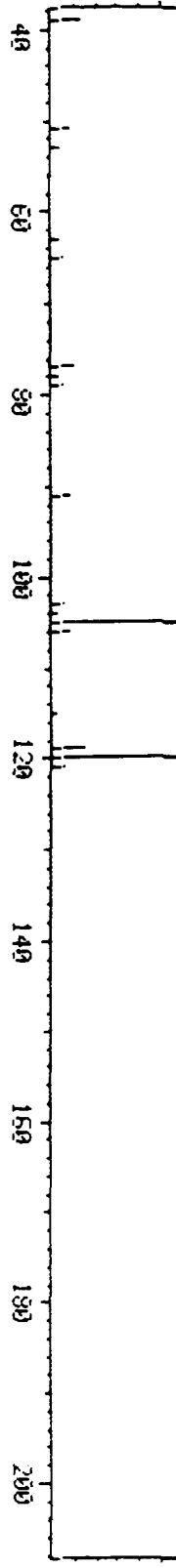
C9.H12
 M UT 120
 B PK 105
 RANK 1
 IN 2527
 PUR 830



C9.H12
 M UT 120
 B PK 105
 RANK 2
 IN 2523
 PUR 865



C9.H12
 M UT 120
 B PK 105
 RANK 3
 IN 2526
 PUR 862



LIBRARY SEARCH
 12/29/89 20:10:00 + 14:56
 SAMPLE: 100ML CC#309686 EPA#B2018 CASE#18765 ON #13
 ENHANCED (S 158 2N 0T)

COMPUchem LABS
 DATA: CSP09686813 #1195
 BASE M/E: 105
 RIC: 120063.

1073
 SAMPLE

C9.H12

M WT 1073
 B PK 105
 RANK 1
 IN 2528
 PUR 687

BENZENE, 1-ETHYL-2-METHYL-

CAS# 611-14-3

C9.H12

M WT 1073
 B PK 105
 RANK 2
 IN 2529
 PUR 679

BENZENE, 1-ETHYL-3-METHYL-

CAS# 620-14-4

C9.H12

M WT 1073
 B PK 105
 RANK 3
 IN 2524
 PUR 676

BENZENE, (1-METHYLETHYL)-

CAS# 98-82-8

M/E

50

100

150

200

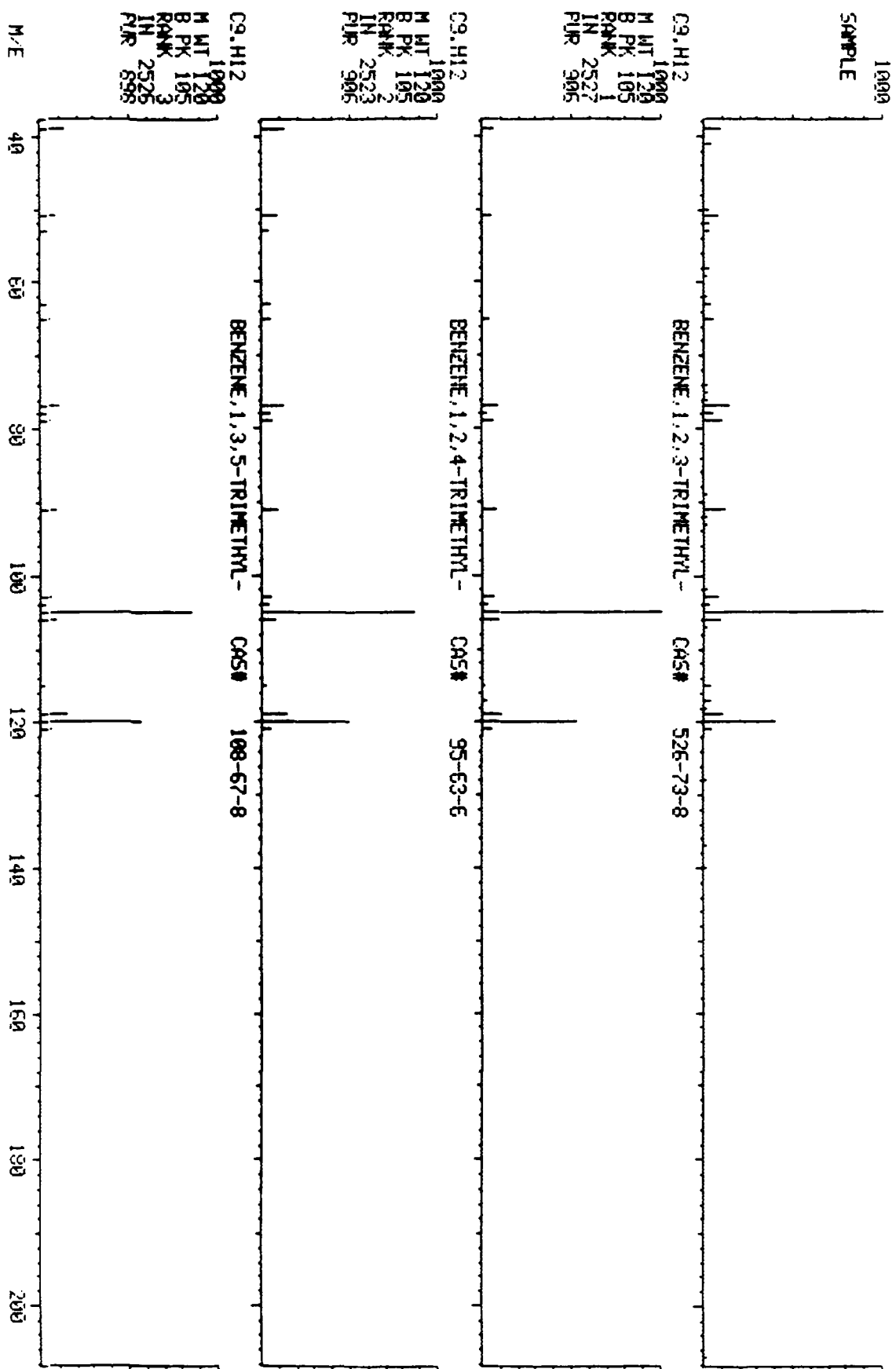
250

LIBRARY SEARCH
 12/29/89 20:10:00 + 15:18
 SAMPLE: 100UL CC#309586 EPAWB2018 CASE#18756 ON #13
 ENHANCED (S 158 2H 0T)

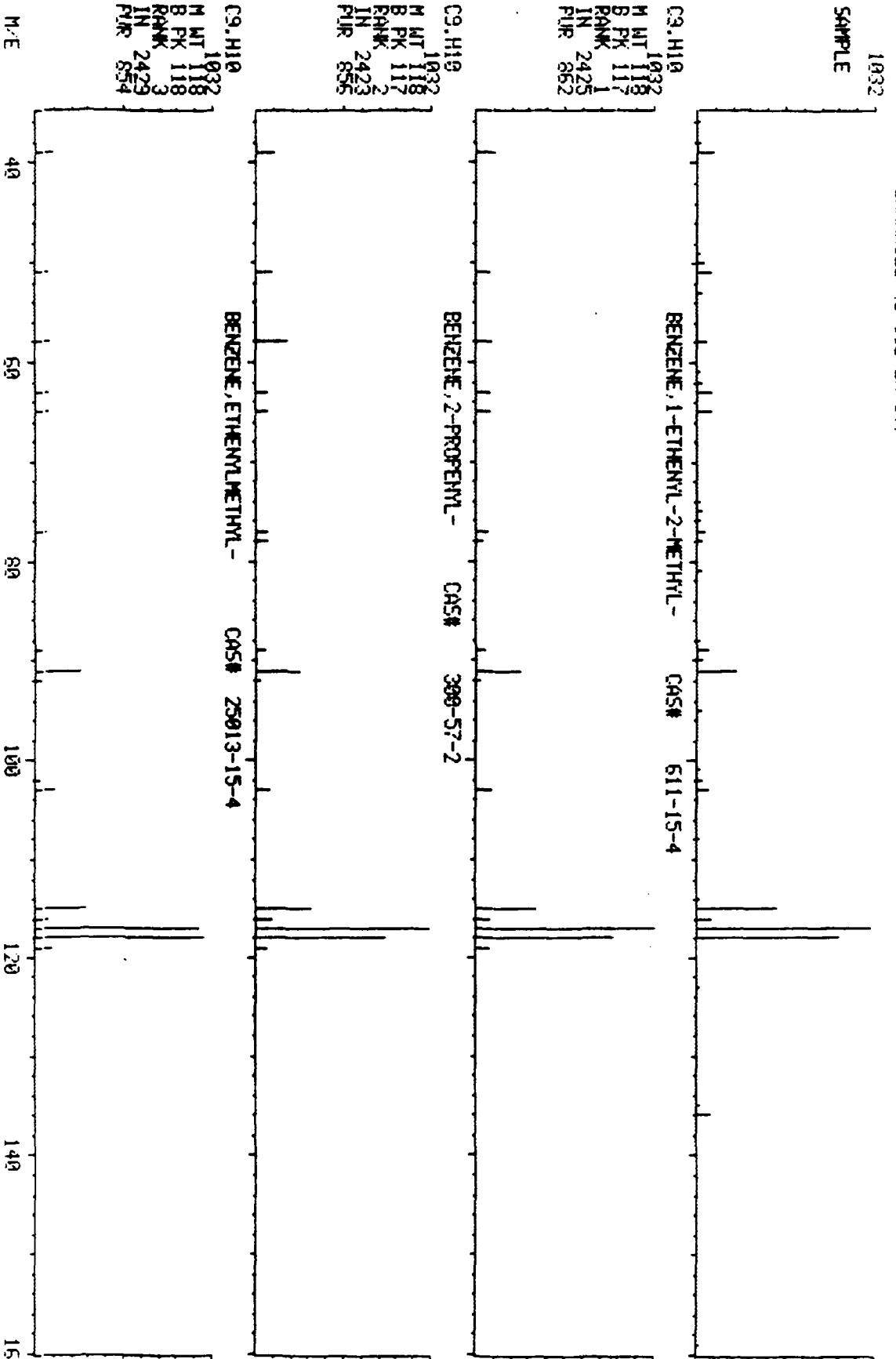
COMPUCHEM LABS

DATA: C5R0368813 #1224

BASE M/E: 105
 RIC: 217599.



COMPUchem LABS
 LIBRARY SEARCH
 12/29/89 20:10:00 + 15:28
 SAMPLE: 100UL CC#309686 EPA#B2018 CASE#19765 ON #13
 ENHANCED (5 158 2H 0T)
 DATA: C5R09686B13 #1237
 BASE M/E: 117
 RIC: 26943.



AB INSTRUCTIONS:

RECEIPT DATE

CASE: 18765 56 *RE 11/2/90*

DUE DATE

VDA
GC/MS WORKSHEET

COMPUCHENG: 3096862

RI 3 R2I 3 D (:1)
R3I 3 R4I 3 D2I 3 (:1)

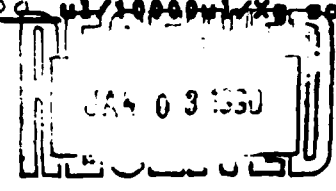
MED. SOLID. EPA 80W 2/87

Sample Prep Code---156
Instrument Code---414
Compound List-----495
Surrogate Std-----394
Internal Std-----036

SAMPLE ID: B2013 RE *RES 250* Dry Wt. Factor 1.14 % Moisture 12

GC/MS ANALYSIS

Amount Purged: [] 10ul ~~9X5~~ soil or [] Dilution 100 ul / 10000 ul ~~9X5~~ soil
Internal Standard Volume Added 5 ul
Surrogate Standard Volume Added - ul
FB Filename BF891229A13 Disk ()
Blank Filename CNO11532B13 Disk ()
Standard Filename CS971229A13 Disk ()
Sample Filename CSR09686B13 Disk ()



ANALYST(S): Injection 1452 Neptune Work-up 1452 Neptune

GC/MS REVIEW

CONDITION CODE

Box containing handwritten codes: RES, EJA

Entry Codes OK, EA, ES, SM, JS, SL, SH, JA, DA

Non-Entry Codes IF, IL, IH, SW, CT, CS, PC, NR, LA, DI, CO, RN, DU, SI, SF, LP, BB, OT, VC, FO, NS

Disposition: [/] Complete

Extraneous Peak Search Results:

of Peaks Found: 10

[] Reprep neat required

Quality Assurance Notice(s):

Notices Required 1



[] Reprep using 9

[] Dilute (:1)

COMMENTS:

GC/MS Review OK Date 1/2/90 Auditor DWagner Date 1/3/90

REPORT INTEGRATION

Final Reportable Package(s): CSR-813 / GRD-C19 Total # of Injections: 6

QA COMMENTS:

FINAL REVIEW:

Initials _____ Date ____/____/____

Initials _____ Date ____/____/____

AC1010 (4/89)

VOLATILE PREPARATION WORKSHEET

QUEUE # 19 PREP CODE -156 ASSIGNED TO MARSHAD Jashi DATE 12/29/89

SAMPLE NUMBER	CASE NUMBER	SDG	QC SAMPLE		SAMPLE WEIGHT (G) VOLUME (ML)	SAMPLE ID	COMMENTS
			TYPE	ORIGINAL			
309686R	18756	0007			4.0g	B2018	
311271	↓	↓	SS	309686	4.0g	B2018MS	
311272	↓	↓	SS	309686	4.0g	B2018MSD	
311273	↓	↓	BS	309686	10.0ml		
311532			B1		10.0ml	I6	
311533			B2		10.0ml	I7	
			B3				
			B4				
			B5				

SURROGATE # 394 LOT # 30837 MANUAL OPERATOR 742 , 267
 AMOUNT 0.5ml
 RELINQUISHED BY ln Jan DATE 12/29/89 RECEIVED BY 1556 DATE 12/29/89

VOLATILE - MEDIUM LEVEL SOLID

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	409	54500	50.0		1200
221	50	CHLOROMETHANE				BDL	1200
231	62	VINYL CHLORIDE				BDL	1200
220	94	BROMOMETHANE				BDL	1200
209	64	CHLOROETHANE				BDL	1200
216	96	1,1-DICHLOROETHENE				BDL	620
254	76	CARBON DISULFIDE				BDL	620
252	43	ACETONE (2-PROPANONE)			1.4	30-100	1200
248	114 I	1,4-DIFLUOROBENZENE (IS)	540	190000	50.0		620
222	84	METHYLENE CHLORIDE			3.8	50-400 D	620
226	96	TRANS-1,2-DICHLOROETHENE				BDL	620
214	63	1,1-DICHLOROETHANE				BDL	620
257	43	VINYL ACETATE				BDL	1200
237	96	CIS-1,2-DICHLOROETHENE				BDL	620
253	72	2-BUTANONE				BDL	1200
211	83	CHLOROFORM				BDL	620
227	97	1,1,1-TRICHLOROETHANE				BDL	620
206	117	CARBON TETRACHLORIDE				BDL	620
203	78	BENZENE			15.2	200-1000 D	620
215	62	1,2-DICHLOROETHANE				BDL	620
270	117 I	D5-CHLOROENZENE (IS)	914	270000	50.0		620
229	130	TRICHLOROETHENE				BDL	620
217	63	1,2-DICHLOROPROPANE				BDL	620
212	83	BROMODICHLOROMETHANE				BDL	620
218	75	CIS-1,3-DICHLOROPROPENE				BDL	620
256	43	4-METHYL-2-PENTANONE			2.0	BDL-250	1200
225	92	TOLUENE			112.0	1000-14000 D	620
250	75	TRANS-1,3-DICHLOROPROPENE				BDL	620
228	97	1,1,2-TRICHLOROETHANE				BDL	620
224	164	TETRACHLOROETHENE				BDL	620
255	43	2-HEXANONE			16.8	200-2100 D	1200
208	129	DIBROMOCHLOROMETHANE				BDL	620
207	112	CHLOROBENZENE				BDL	620
219	106	ETHYLBENZENE			52.0	700-6500 D	620
330	106	M, P-XYLENE			86.6	1200-11000 D	620
239	106	O-XYLENE			60.4	800-7600 D	620
251	104	STYRENE			19.6	200-2400 D	620
205	173	BROMOFORM				BDL	620
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	620
258	65 S	D4-1,2-DICHLOROETHANE NA#57			56.4	113. %	
247	95 S	BROMOFLUOROBENZENE			52.3	105. %	
233	98 S	D8-TOLUENE NA#59			57.3	115. %	
289	106	XYLENES (TOTAL)			147.0	2000-10000 D	620

CORRECTED/REVIEWED BY *CK St...*
(GC/MS DATA REVIEWER)

DATE 12-98

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
299	96	1,2-DICHLOROETHENE (TOTAL)				BDL	620
CHECKSUMS:							
		3979.	1863	514500.	832.8	64743	

CORRECTED/REVIEWED BY *CK Stuber*
(GC/MS DATA REVIEWER)

DATE 1-2-90

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P	F
40	258	D4-1,2-DICHLOROETHANE NA#57	56.4	50.0	113.	70-121	X	
41	247	BROMOFLUOROBENZENE	52.3	50.0	105.	74-121	X	
42	233	D8-TOLUENE NA#59	57.3	50.0	115.	81-117	X	

* ADVISORY SURROGATE ONLY

++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

CORRECTION FACTOR CALCULATION:

$$\frac{4.0 \text{ G}}{\text{WET WEIGHT OF SAMPLE (G)}} \times \text{LIBRARY ADJUSTMENT} \times \text{DRY WEIGHT FACTOR} \times \frac{10000 \text{ UL}}{\text{UL USED}} =$$

$$\frac{4.0 \text{ G}}{4.00 \text{ (G)}} \times 1.25 \times \frac{1.14}{1.00} \times \frac{10000 \text{ UL}}{100. \text{ UL}} = 125.000$$

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:

$$\frac{100. \text{ UL}}{\text{UL USED}} =$$

$$\frac{100. \text{ UL}}{100. \text{ UL}} = 1.0$$



VERSION 8

CORRECTED/REVIEWED BY *Ok ftd r*
(GC/MS DATA REVIEWER)

DATE) 2.90

LABORATORY NOTICE

CompuChem ID #309686 (Client ID#B2018)

Case# 18756

The volatile fraction of this sample was analyzed at 1.0 gram as a Low Level Solid. The 1.0 gram analysis indicated a Medium Level Solid preparation was necessary to bring concentrations of the target compounds Toluene, 2-Hexanone, M,P-Xylene, and O-Xylene within the analytical range. The analysis of 100 ul of the Medium Level extract should have resulted in quantitative report concentrations in the upper half of the analytical range. However, results varied between the two analyses as shown in the following table of reported amounts:

	Low Level Soil		Medium Level Soil	
Acetone	370	ug/kg	BDL*	
Methylene Chloride	69	ug/kg	540	ug/kg
2-Butanone	260	ug/kg	BDL	
Benzene	140	ug/kg	2200	ug/kg
4-Methyl-2-Pentanone	71	ug/kg	BDL	
Toluene	1600	ug/kg	16000	ug/kg
2-Hexanone	1600	ug/kg	2400	ug/kg
Ethylbenzene	980	ug/kg	7400	ug/kg
M,P-Xylene	1900	ug/kg	12000	ug/kg
O-Xylene	1400	ug/kg	8600	ug/kg
Styrene	540	ug/kg	2800	ug/kg

* = Below Detection Limits

We are attributing these variations to an inhomogeneous soil matrix as well as to differences in the Low Level Soil Purge and Trap analysis and the Medium Level Soil methanol extraction procedure. We are reporting these data with reference to this notice.

Owen K. Studt
Data Review Specialist
15 January, 1990

QA Approval #1032
Linda Fowler
Sr. Quality Assurance Specialist
January 16, 1990

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309688
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009688C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/22/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NO.	COMPOUND	(ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	11	U
74-83-9	-----Bromomethane	11	U
75-01-4	-----Vinyl Chloride	11	U
75-00-3	-----Chloroethane	11	U
75-09-2	-----Methylene Chloride	33	B
67-64-1	-----Acetone	37	B
75-15-0	-----Carbon Disulfide	6	U
75-35-4	-----1,1-Dichloroethene	6	U
75-34-3	-----1,1-Dichloroethane	6	U
540-59-0	-----1,2-Dichloroethene (total)	6	U
67-66-3	-----Chloroform	6	U
107-06-2	-----1,2-Dichloroethane	6	U
78-93-3	-----2-Butanone	26	U
71-55-6	-----1,1,1-Trichloroethane	6	U
56-23-5	-----Carbon Tetrachloride	6	U
108-05-4	-----Vinyl Acetate	11	U
75-27-4	-----Bromodichloromethane	6	U
78-87-5	-----1,2-Dichloropropane	6	U
10061-01-5	-----cis-1,3-Dichloropropene	6	U
79-01-6	-----Trichloroethene	6	U
124-48-1	-----Dibromochloromethane	6	U
79-00-5	-----1,1,2-Trichloroethane	6	U
71-43-2	-----Benzene	6	U
10061-02-6	-----Trans-1,3-Dichloropropene	6	U
75-25-2	-----Bromoform	6	U
108-10-1	-----4-Methyl-2-Pentanone	11	U
591-78-6	-----2-Hexanone	13	U
127-18-4	-----Tetrachloroethene	6	U
79-34-5	-----1,1,2,2-Tetrachloroethane	6	U
108-88-3	-----Toluene	1	J
108-90-7	-----Chlorobenzene	6	U
100-41-4	-----Ethylbenzene	6	U
100-42-5	-----Styrene	6	U
1330-20-7	-----Total Xylenes	2	J

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202A

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309688
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009688C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 12 Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 4

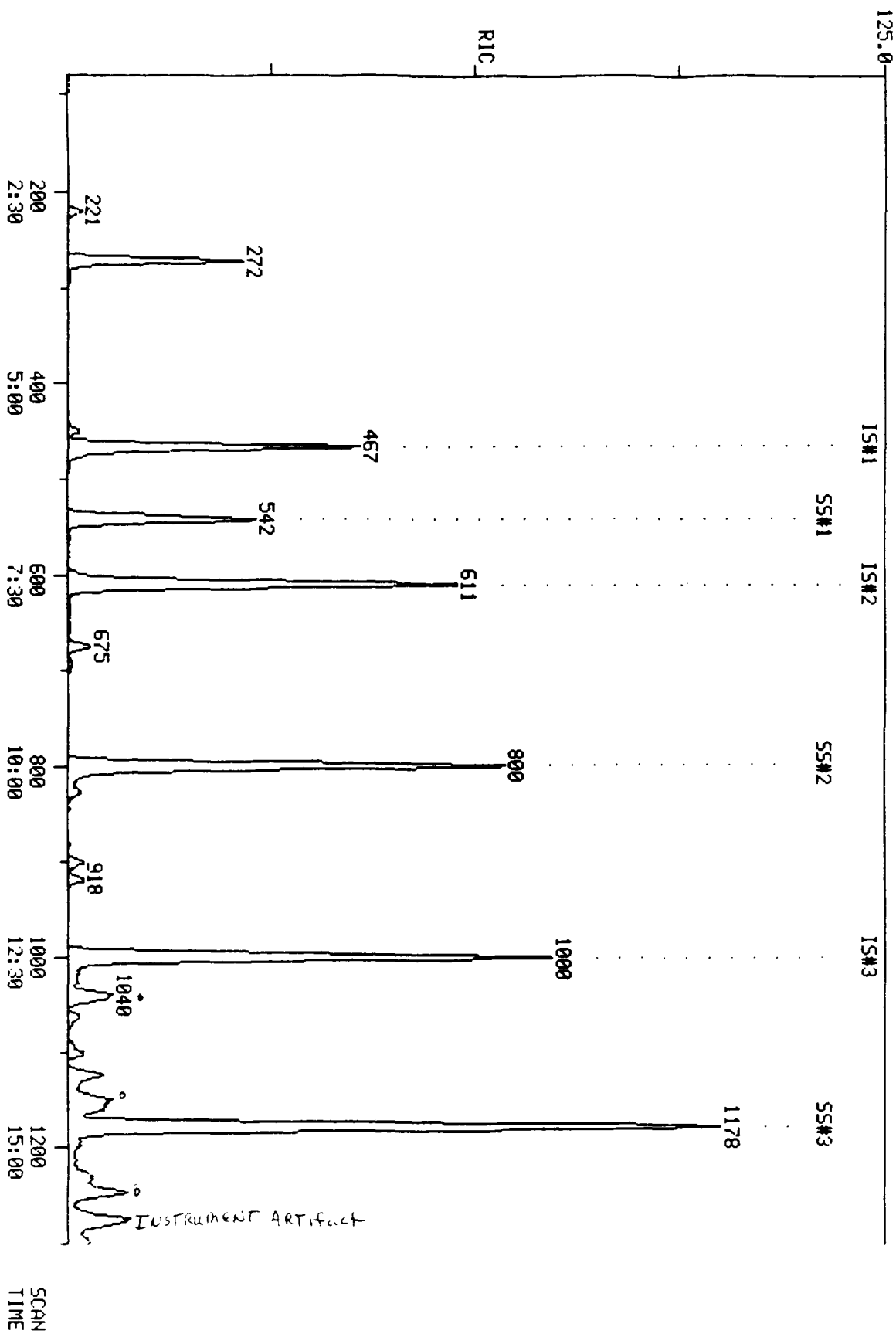
CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	13.00	6.8	J
2.	UNKNOWN	14.37	5.7	J
3.	UNKNOWN	15.57	6.8	J
4.	INSTRUMENT ARTIFACT	15.92	6.8	J

RIC
12/22/89 6:56:00
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
COND5.:

COMPUCHEM LABS
COMPUCHEM DATA: GH009588C19 SCANS 80 TO 1300

119840.



QUANTITATION REPORT FILE: GH009688C19
 DATA: GH009688C19.TI
 12/22/89 6:56:00
 SAMPLE: 5G CC#309683 CASE#18756.7 EPA#B202A ON#19
 CONDS.:
 SUBMITTED BY: 19 ANALYST: 1422

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
 RESP. FAC. FROM LIBRARY ENTRY -

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1
2	221 CHLOROMETHANE <74-87-3> RO#2
3	231 VINYL CHLORIDE <75-01-4> RO#3
4	220 BROMOMETHANE <78-83-9> RO#4
5	209 CHLOROETHANE <75-00-3> RO#5
6	216 1,1-DICHLOROETHENE <75-35-4> RO#8
7	254 CARBON DISULFIDE <75-15-0> RO#9
8	252 ACETONE (2-PROPANONE) <67-64-1> RO#13
9	*248 1,4-DIFLUOROBENZENE (IS) <540-36-3> RO#14
10	222 METHYLENE CHLORIDE <75-09-2> RO#16
11	226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17
12	214 1,1-DICHLOROETHANE <75-34-3> RO#19
13	257 VINYL ACETATE <108-05-4> RO#20
14	237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21
15	253 2-BUTANONE <78-93-3> RO#22
16	211 CHLOROFORM <67-66-2> RO#23
17	227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24
18	206 CARBON TETRACHLORIDE <56-23-5> RO#25
19	203 BENZENE <71-43-2> RO#26
20	215 1,2-DICHLOROETHANE <107-06-2> RO#27
21	*270 D5-CHLOROBENZENE (IS) RO#29
22	229 TRICHLOROETHENE <79-01-6> RO#30
23	217 1,2-DICHLOROPROPANE <78-87-5> RO#31
24	212 BROMODICHLOROMETHANE <75-27-4> RO#33
25	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35
26	256 4-METHYL-2-PENTANONE <108-01-1> RO#36
27	225 TOLUENE <108-88-3> RO#37
28	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38
29	228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39
30	224 TETRACHLOROETHENE <127-18-4> RO#41
31	255 2-HEXANONE <591-78-6> RO#42
32	208 DIBROMOCHLOROMETHANE <124-48-1> RO#43
33	207 CHLOROBENZENE <108-90-7> RO#45
34	219 ETHYLBENZENE <100-41-4> RO#47
35	330 M,P-XYLENE <133-02-7> RO#48
36	239 O-XYLENE <133-02-7> RO#49
37	251 STYRENE <100-42-5> RO#50
38	205 BROMOFORM <75-25-2> RO#51
39	223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54
40	*259 D4-1,2-DICHLOROETHANE RO#57
41	*247 BROMOFLUROBENZENE <460-00-4> RO#58
42	*233 D8-TOLUENE RO#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	467	5:50	1	1.000	A BB	49330.	50.000 UG/KG	12.64
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	222	2:46	1	0.475	A BB	8503.	32.570 UG/KG	8.23 ⁴²
9	114	611	7:38	9	1.000	A BB	196249.	50.000 UG/KG	12.64
10	84	271	3:23	1	0.580	A BB	47558.	29.334 UG/KG	7.41 ⁵⁰
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	451	5:38	1	0.966	A BB	2422.	22.907 UG/KG	5.79 ⁴²
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	NOT FOUND							
20	62	NOT FOUND							
21	117	999	12:29	21	1.000	A BB	206316.	50.000 UG/KG	12.64
22	130	NOT FOUND							
23	63	NOT FOUND							
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	801	10:01	21	0.802	A BB	2425.	2.290 UG/KG	0.58 ^{NO}
27	92	808	10:06	21	0.809	A BB	3891	1.278 UG/KG	0.32 ⁴²
28	75	NOT FOUND							
29	97	NOT FOUND							
30	164	NOT FOUND							
31	43	920	11:30	21	0.921	A VB	8912.	11.064 UG/KG	2.80 ⁴²
32	129	NOT FOUND							
33	112	NOT FOUND							
34	105 ²²	1046	13:04	21	1.047	A BB	1681 4995.	0.642 389 UG/KG	0.60 ^{17DC}
35	105	1046	13:04	21	1.047	A BB	4995.	1.718 UG/KG	0.43 ⁴²
36	106	NOT FOUND							
37	104	NOT FOUND							
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	542	6:46	1	1.161	A BB	99647.	48.404 UG/KG	12.23
41	95	1178	14:43	21	1.179	A BB	162654.	46.080 UG/KG	11.65
42	98	800	10:00	21	0.801	A BB	206033.	47.659 UG/KG	12.04

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:54	0.99	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:06		10.000			50.00		0.952	
3	1:11		10.000			50.00		1.256	
4	1:26		10.000			50.00		1.579	
5	1:34		10.000			50.00		0.796	
6	2:37		5.000			50.00		1.346	
7	2:46		5.000			50.00		3.709	
8	2:48	0.99	10.000	0.05	32.57	50.00	0.172	0.265	0.65
9	7:42	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	3:27	0.98	5.000	0.12	29.33	50.00	0.964	1.643	0.59
11	3:53		5.000			50.00		1.486	
12	4:34		5.000			50.00		2.635	
13	4:52		10.000			50.00		0.458	
14	5:33		5.000			50.00		1.675	

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:42	0.99	10.000	0.10	22.91	50.00	0.049	0.107	0.46
16	6:10		5.000			50.00		3.433	
17	6:18		5.000			50.00		0.814	
18	6:34		5.000			50.00		0.755	
19	6:53		5.000			50.00		0.895	
20	6:58		5.000			50.00		2.523	
21	12:34	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	8:01		5.000			50.00		0.569	
23	8:22		5.000			50.00		0.352	
24	8:55		5.000			50.00		0.701	
25	9:41		5.000			50.00		0.690	
26	10:04	0.99	15.000	0.05	2.29	50.00	0.012	0.257	0.05
27	10:10	0.99	5.000	0.16	1.28	50.00	0.019	0.738	0.03
28	10:42		5.000			50.00		0.342	
29	10:58		5.000			50.00		0.349	
30	11:06		5.000			50.00		0.723	
31	11:33	1.00	15.000	0.06	11.06	50.00	0.043	0.195	0.22
32	11:35		5.000			50.00		0.690	
33	12:37		5.000			50.00		1.092	
34	12:55	1.01	5.000	0.21	2.39	50.00	0.024	0.507	0.05
35	13:10	0.99	5.000	0.21	1.72	50.00	0.024	0.705	0.03
36	13:52		5.000			50.00		0.622	
37	13:54		5.000			50.00		1.099	
38	14:07		5.000			50.00		0.664	
39	15:16		5.000			50.00		0.557	
40	6:52	0.99	5.000	0.23	48.40	50.00	2.020	2.087	0.97
41	14:49	0.99	5.000	0.24	46.08	50.00	0.788	0.855	0.92
42	10:04	0.99	5.000	0.16	47.66	50.00	0.999	1.048	0.95

LIBRARY SEARCH
12/22/89 6:56:00 + 2:45
SAMPLE: 5G CC#309688 CASE#18756.7 EPAN#202A ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

DATA: GH009688C19 # 222

BASE M/E: 43
RIC: 1783.

1092
SAMPLE

C3.H6.0
M.WT 1092
B.PK 43
RANK 1
IN 13
PUR 644

252 ACETONE (2-PROPANONE) <67-64-1> RM#13

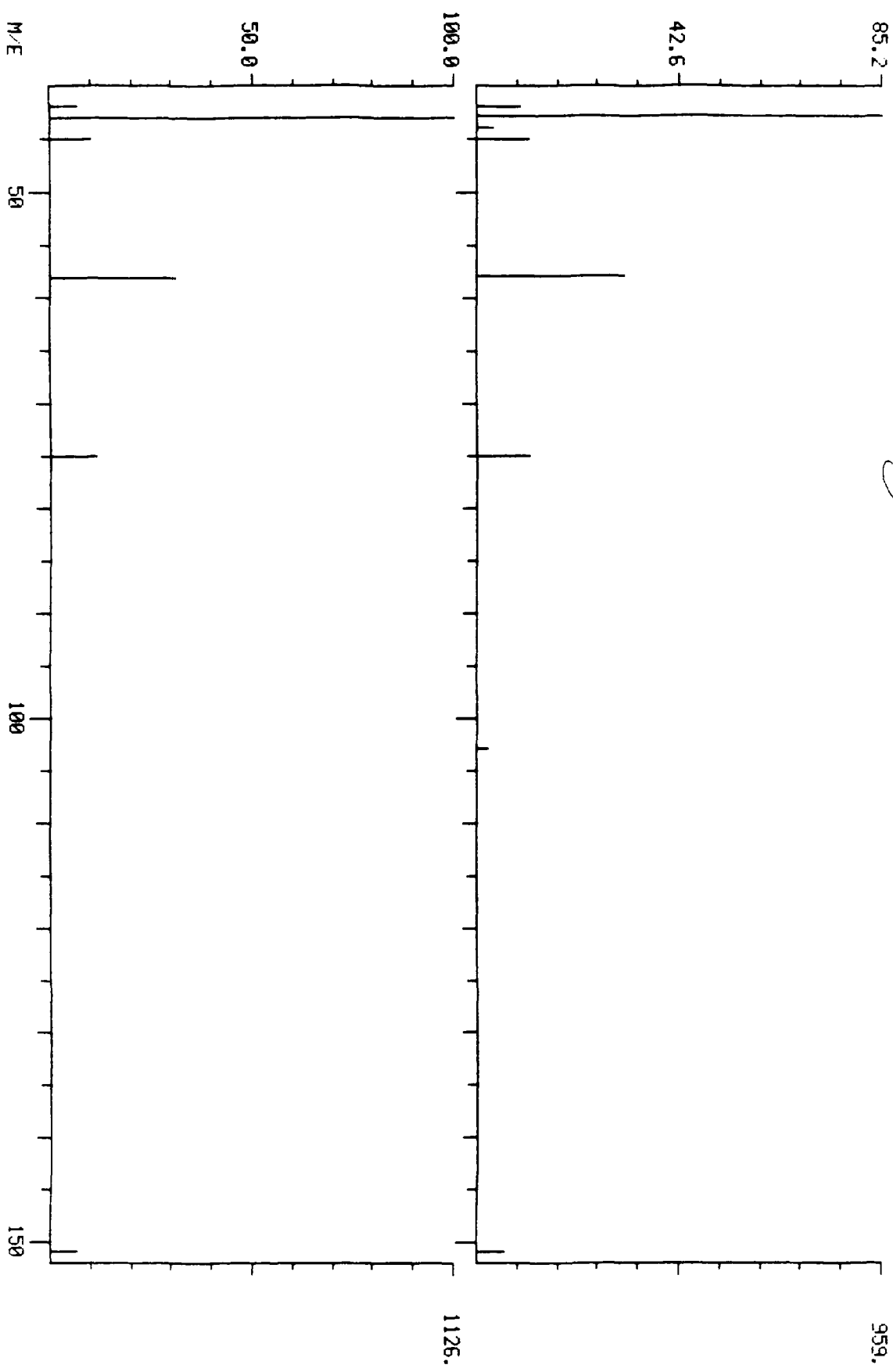
SAMPLE MINUS LIBRARY



DUPL MASS SPECTRUM
12/22/89 6:56:00 + 2:46
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (5 158 2N) (252) ACETONE (2-PROPANONE) (67-64-1) R0#13

COMPUCHEN LABS

DATA: GH009688C19 #222 BASE M/E: 43/ 43
RIC: 1783.7 1869.

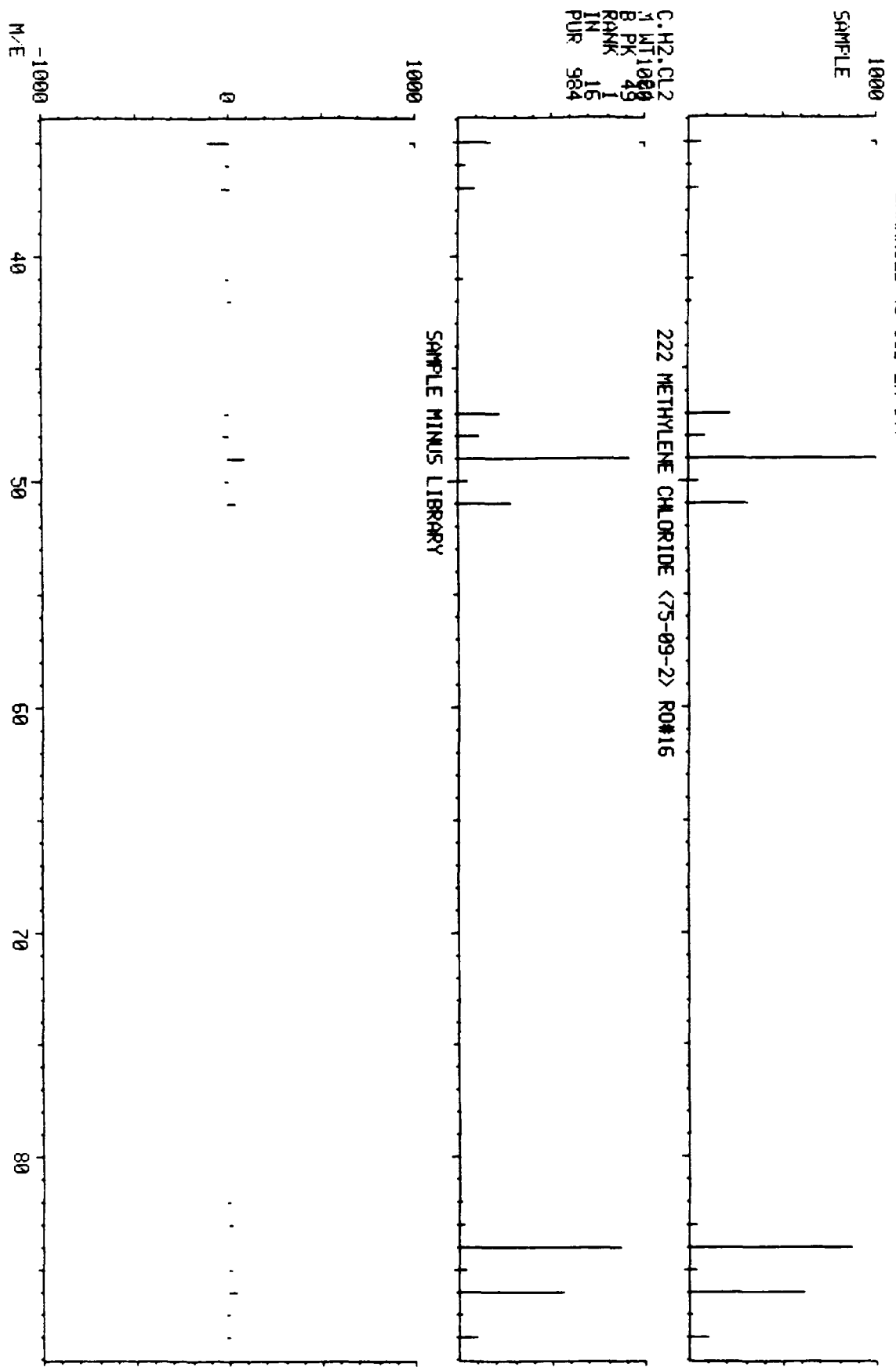


COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 6:56:00 + 3:23
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A QN#19
ENHANCED (5 158 2N 0T)

DATA: GH009688C19 # 271

BASE M/E: 49
RIC: 23519.

C.H2.C12
1 WT 1000
B PK 49
RANK 1
IN 16
PUR 984



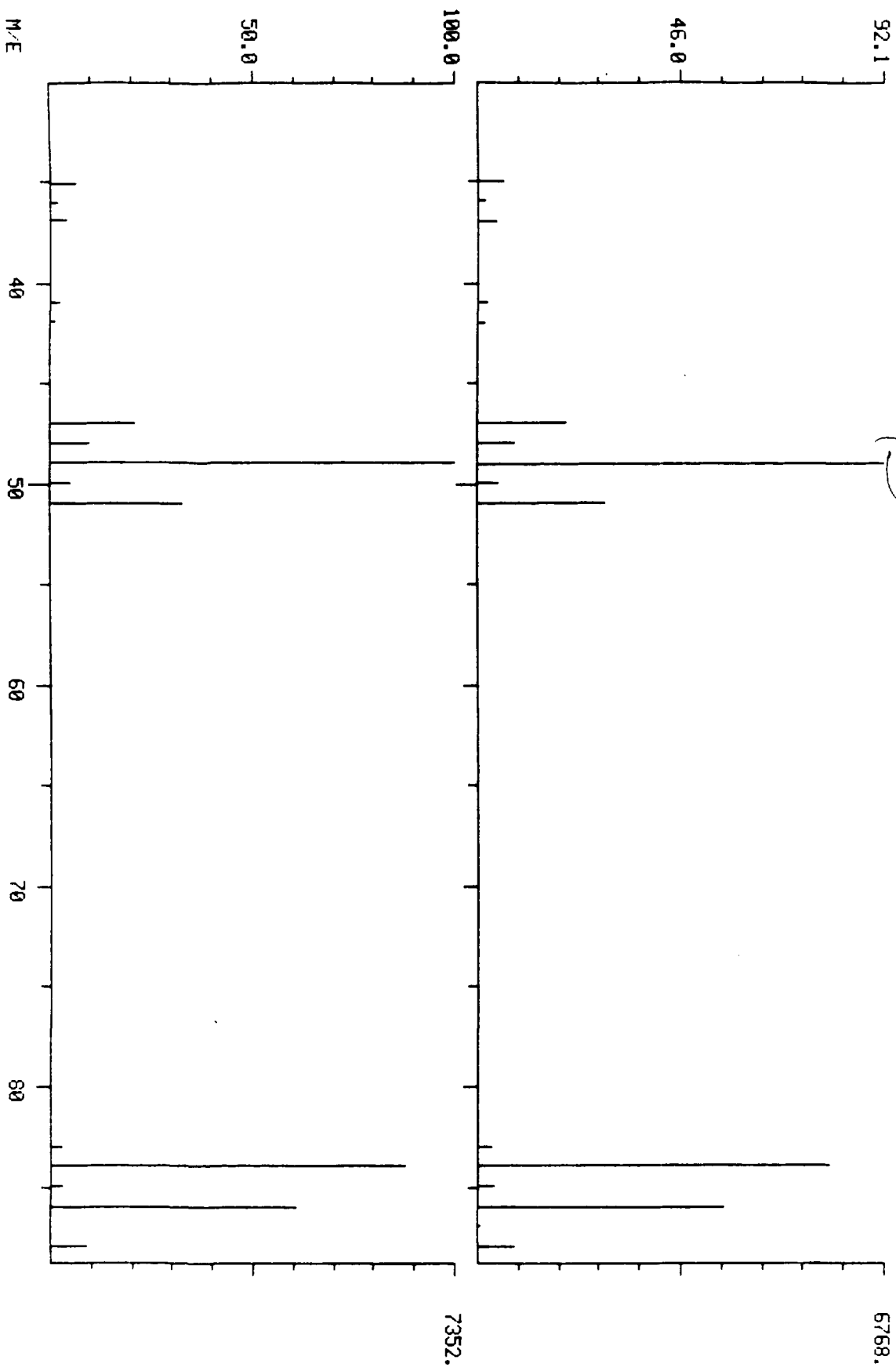
DUAL MASS SPECTRUM
12/22/89 6:56:00 + 3:23
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (S 15B 2N)

COMPUCHEN LABS

DATA: CH009688C19 #271

BASE M/E: 49/ 49
RIC: 23519. / 25439.

222 METHYLENE CHLORIDE <75-09-2> R0#16



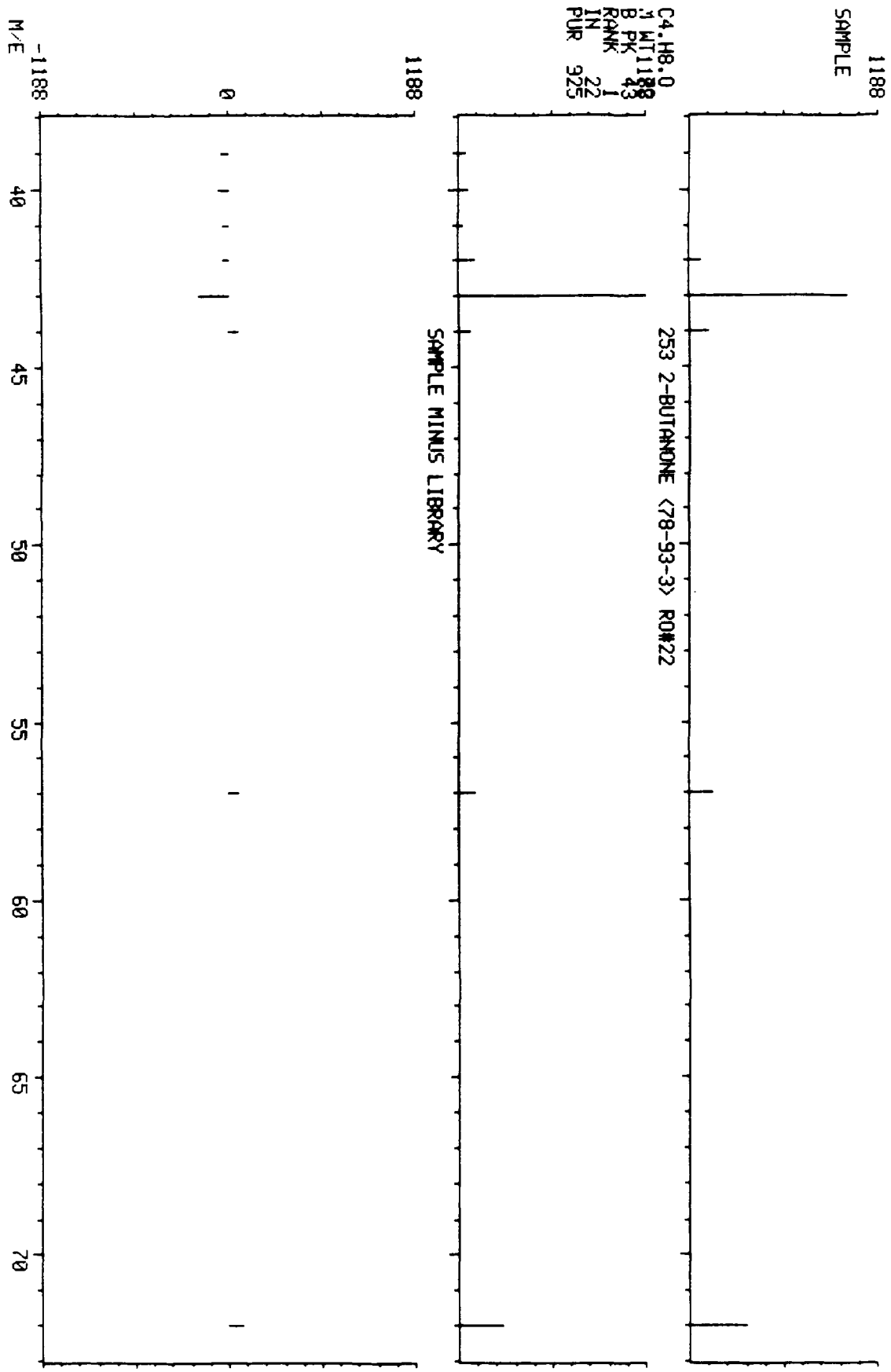
LIBRARY SEARCH
12/22/89 6:55:00 + 5:38
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (S 158 2N 0T)

COMPUCHEM LABS

DATA: GH009688C19 # 451

BASE M/E: 43
RIC: 1463.

C4.H8.0
Y MIT 1188
B PK 43
R RANK 1
IN 22
PUR 925

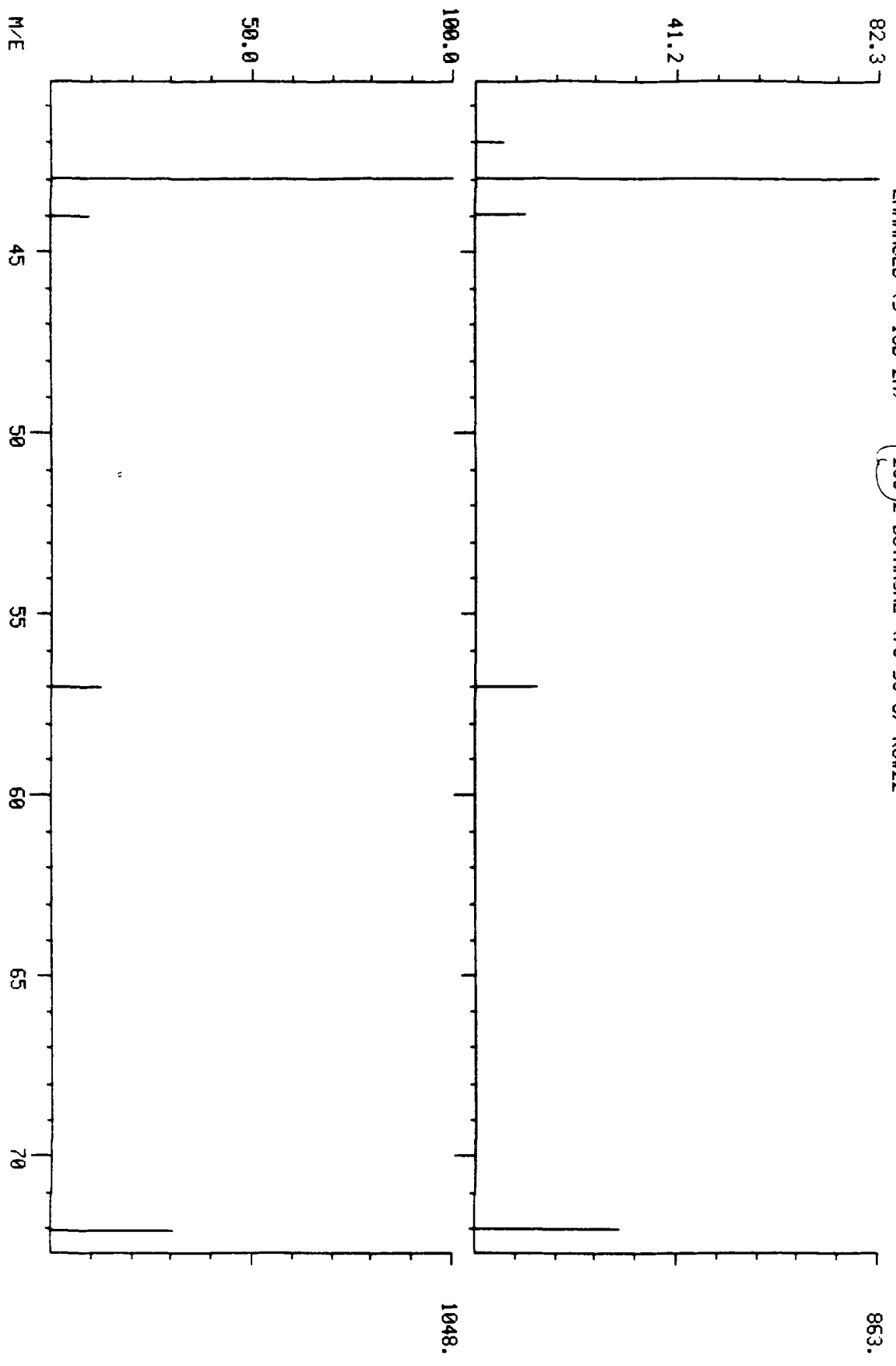


DUAL MASS SPECTRUM
12/22/89 6:56:00 + 5:38
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#8202A ON#19
ENHANCED (5 15B 2N) (253) 2-BUTANONE (78-93-3) R0#22

COMPUchem LABS

DATA: GH009688C19 #451

BASE M/E: 43/ 43
RIC: 1463.7 1591.

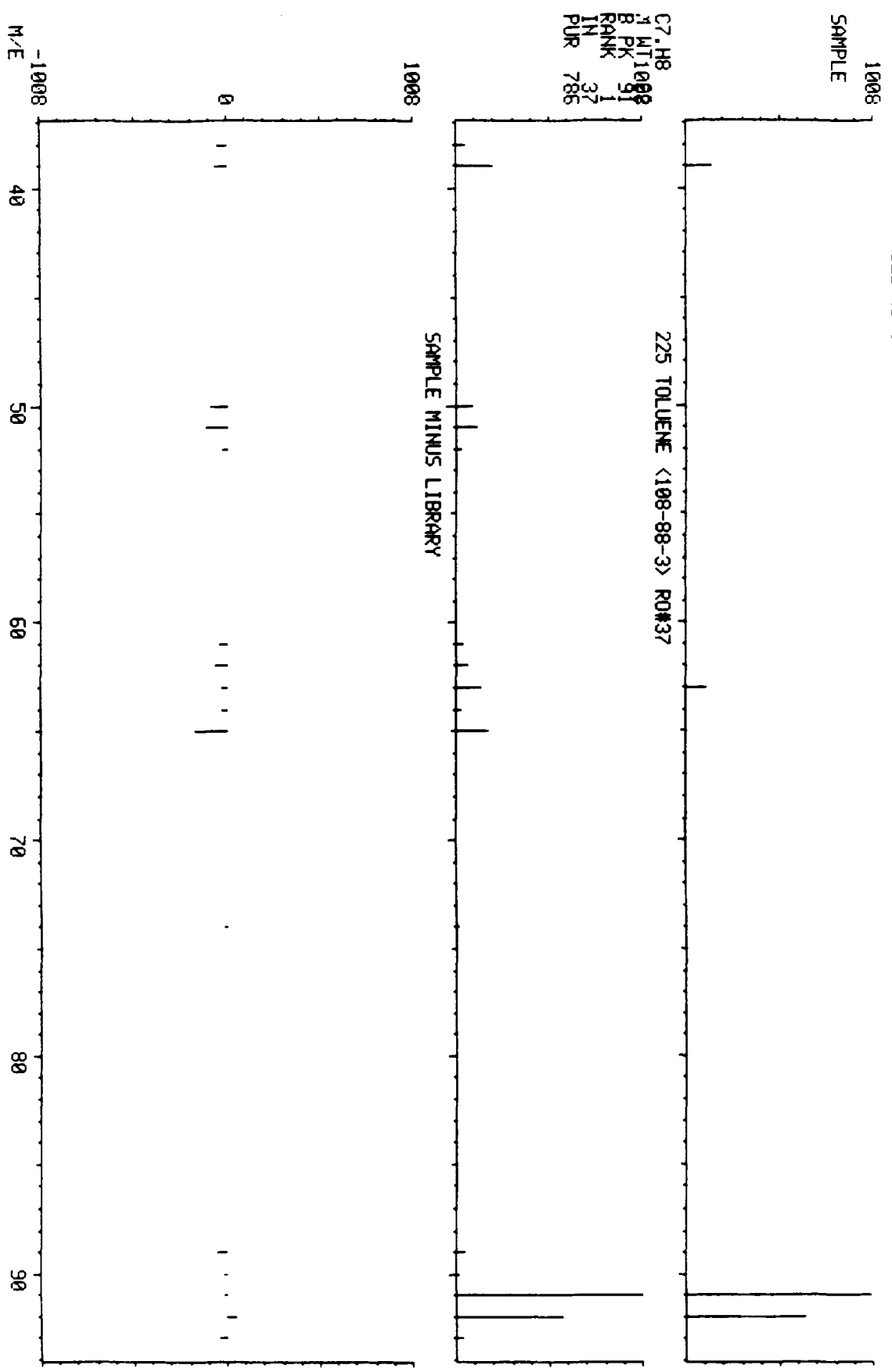


COMPUCEM LAB5
LIBRARY SEARCH
12/22/89 6:56:00 + 10:06
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (5 158 2N 0T)

DATA: GH09689C19 # 808

BASE M/E: 91
RIC: 1167.

C7.H8
1 MT 1008
B PK 91
RANK 37
IN 786
PUR

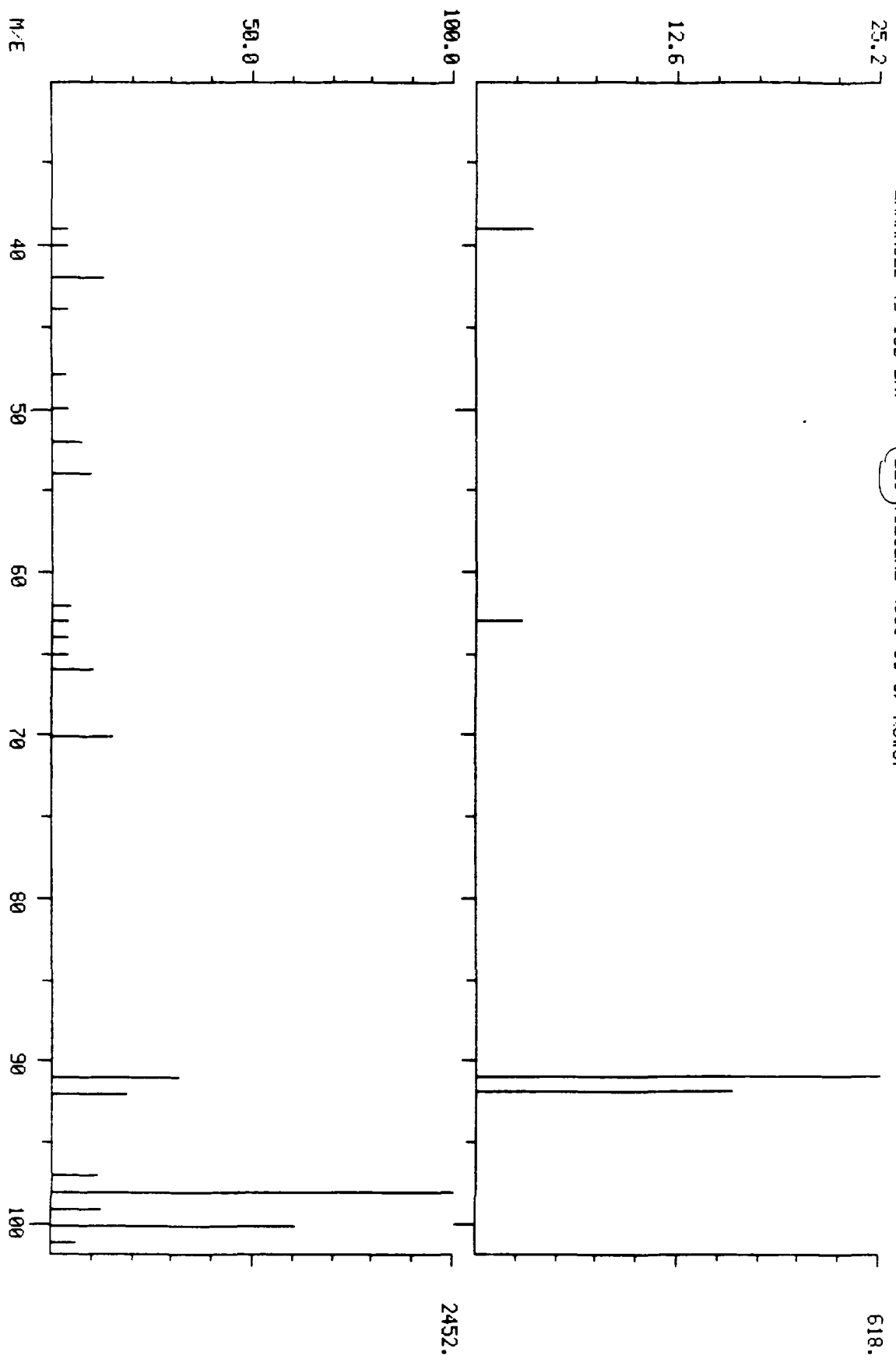


DUAL MASS SPECTRUM
12/22/89 6:56:00 + 10:06
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A DN#19
ENHANCED (5 158 2M) (225) TOLUENE <108-88-3> R0#37

COMPUCHEM LABS

DATA: GH009688C19 #808

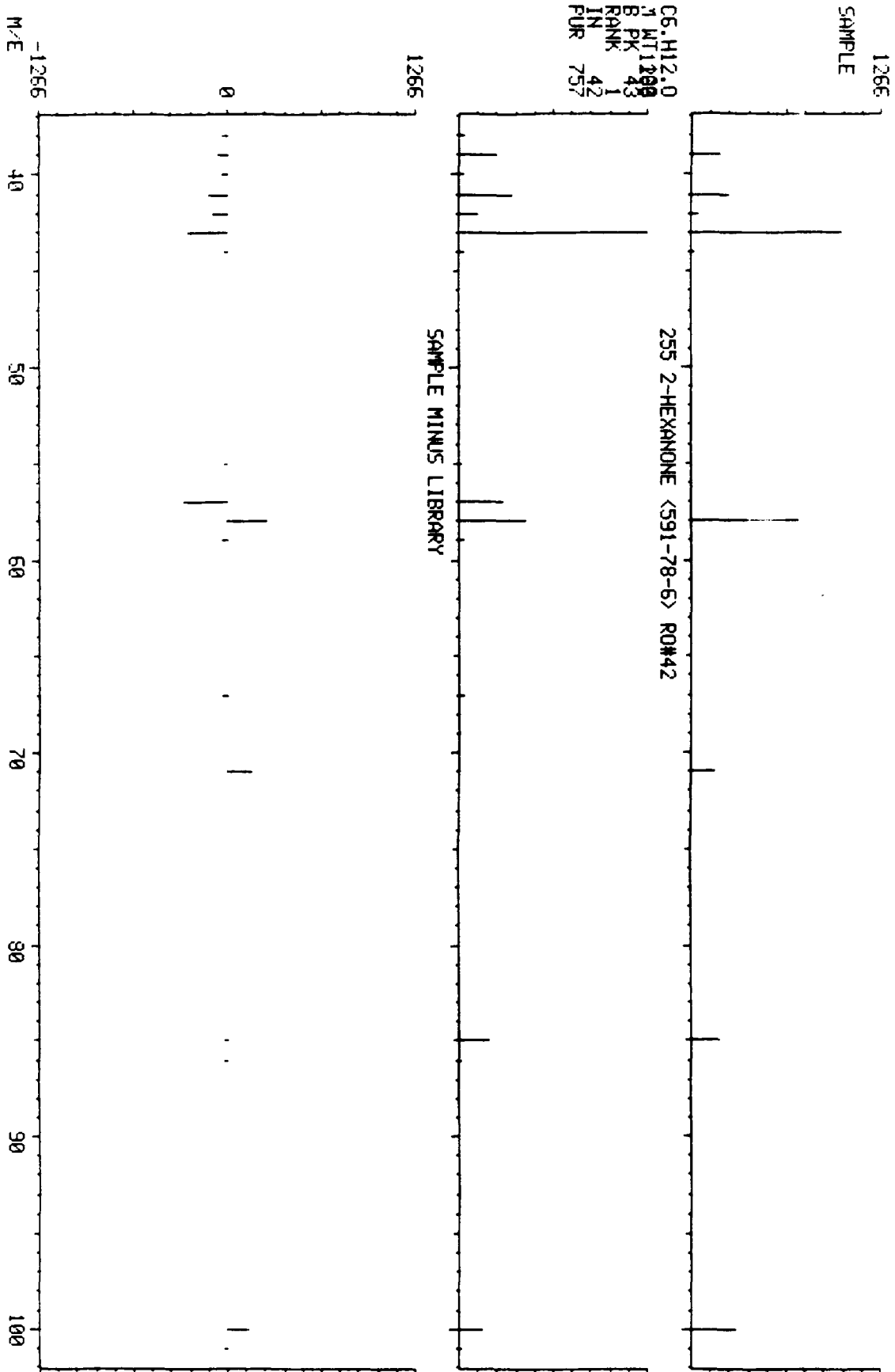
BASE M/E: 91 / 98
RIC: 1167. / 8103.



COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 6:56:00 + 11:30
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (5 158 2H 0T)

DATA: GH009688C19 # 920
BASE M/E: 43
RIC: 1721.

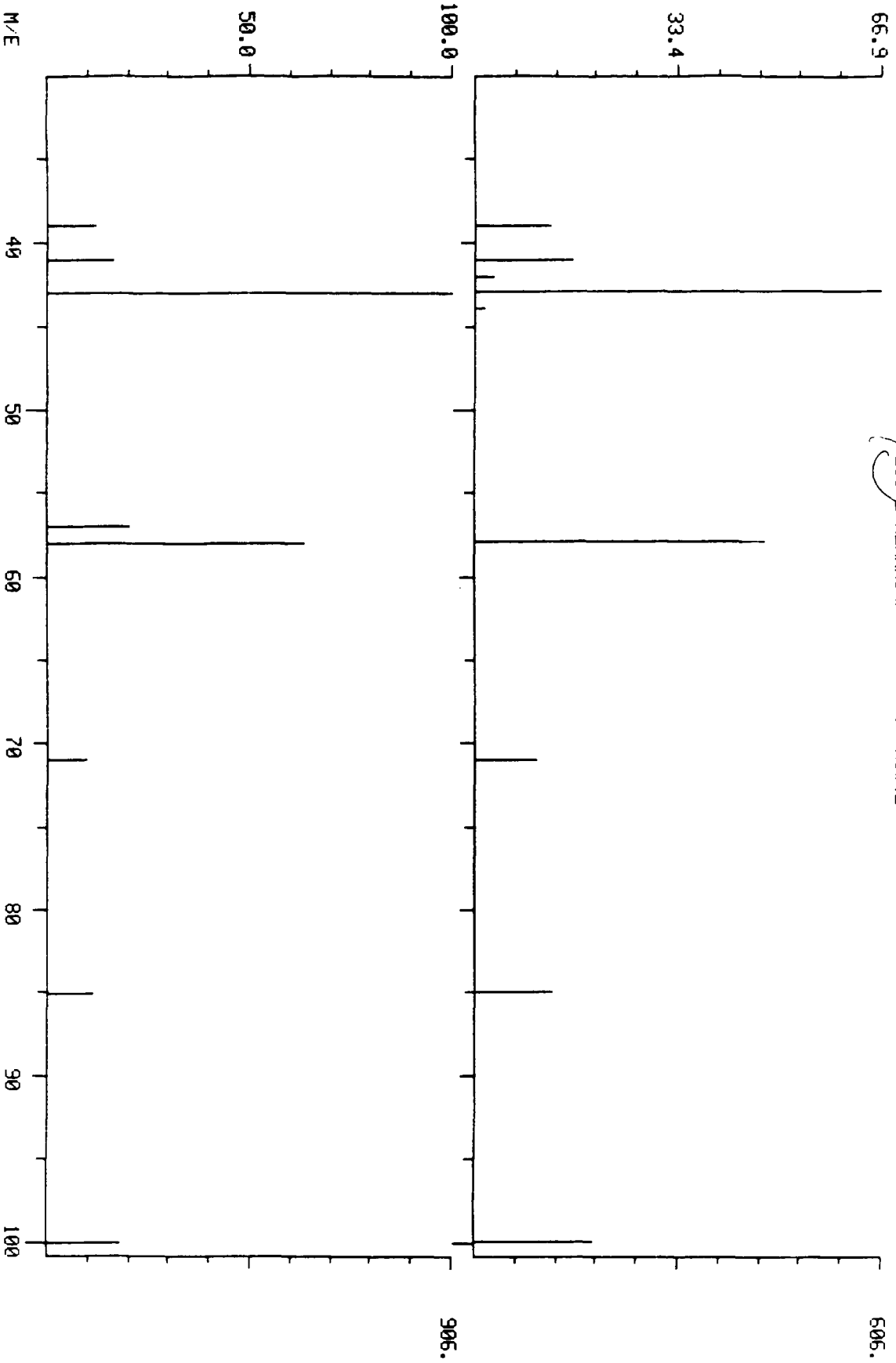
CG.H12.0
1 WT 1298
B PK 43
RANK 1
IN 42
PUR 757



DUAL MASS SPECTRUM
12/22/89 6:56:00 + 11:30
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (5 15B 2N) 255 2-HEXANONE (591-78-6) R0#42

COMPUCHEM LABS

DATA: CH009688C19 #920 BASE M/E: 43/ 43
RIC: 1721. / 2267.

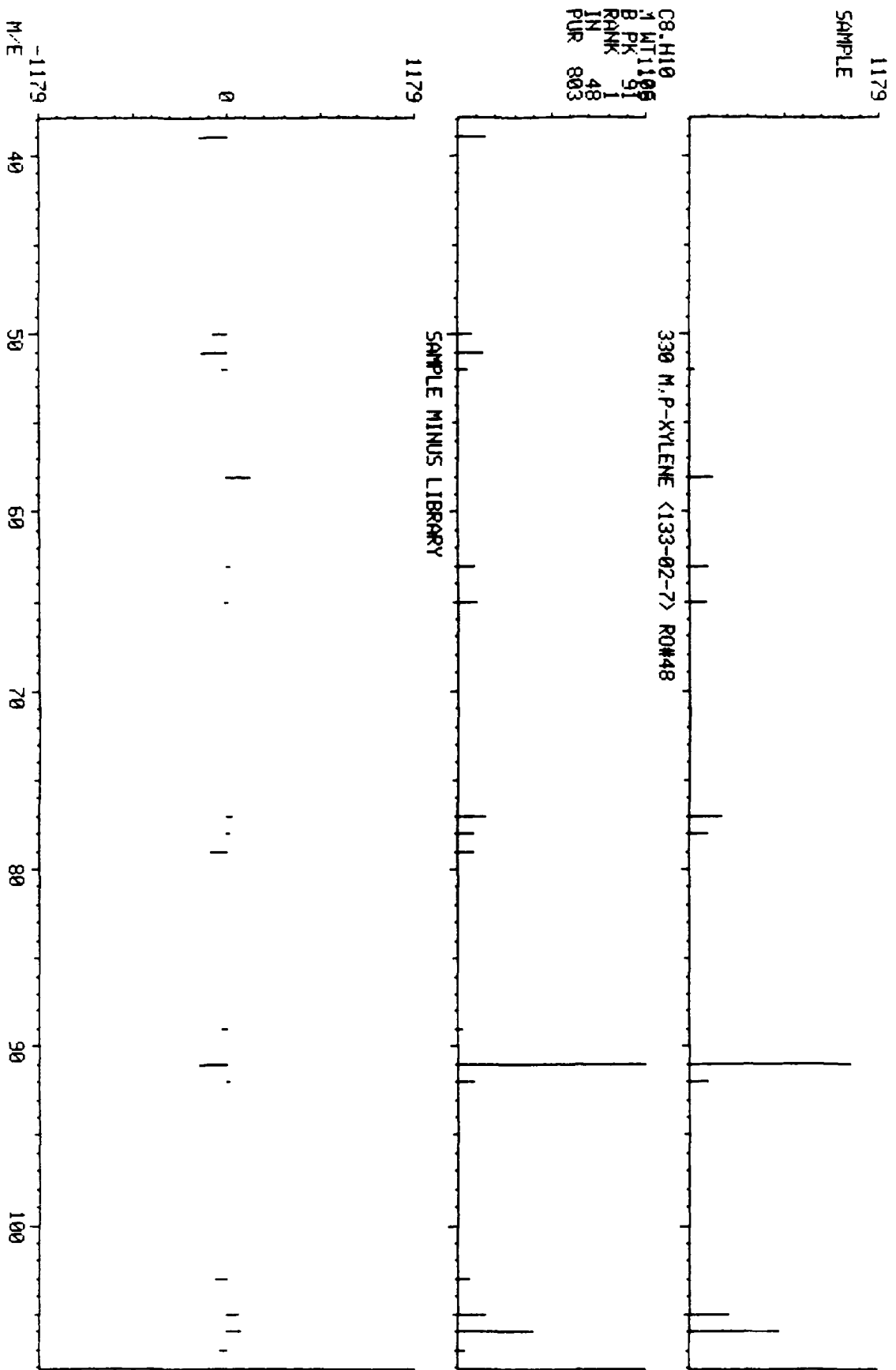


LIBRARY SEARCH
12/22/89 6:56:00 + 13:04
SAMPLE: 5G.CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (S 158 ZN 01)

COMPUCHEM LABS

DATA: GH009688C19 #1046

BASE M/E: 91
RIC: 2363.

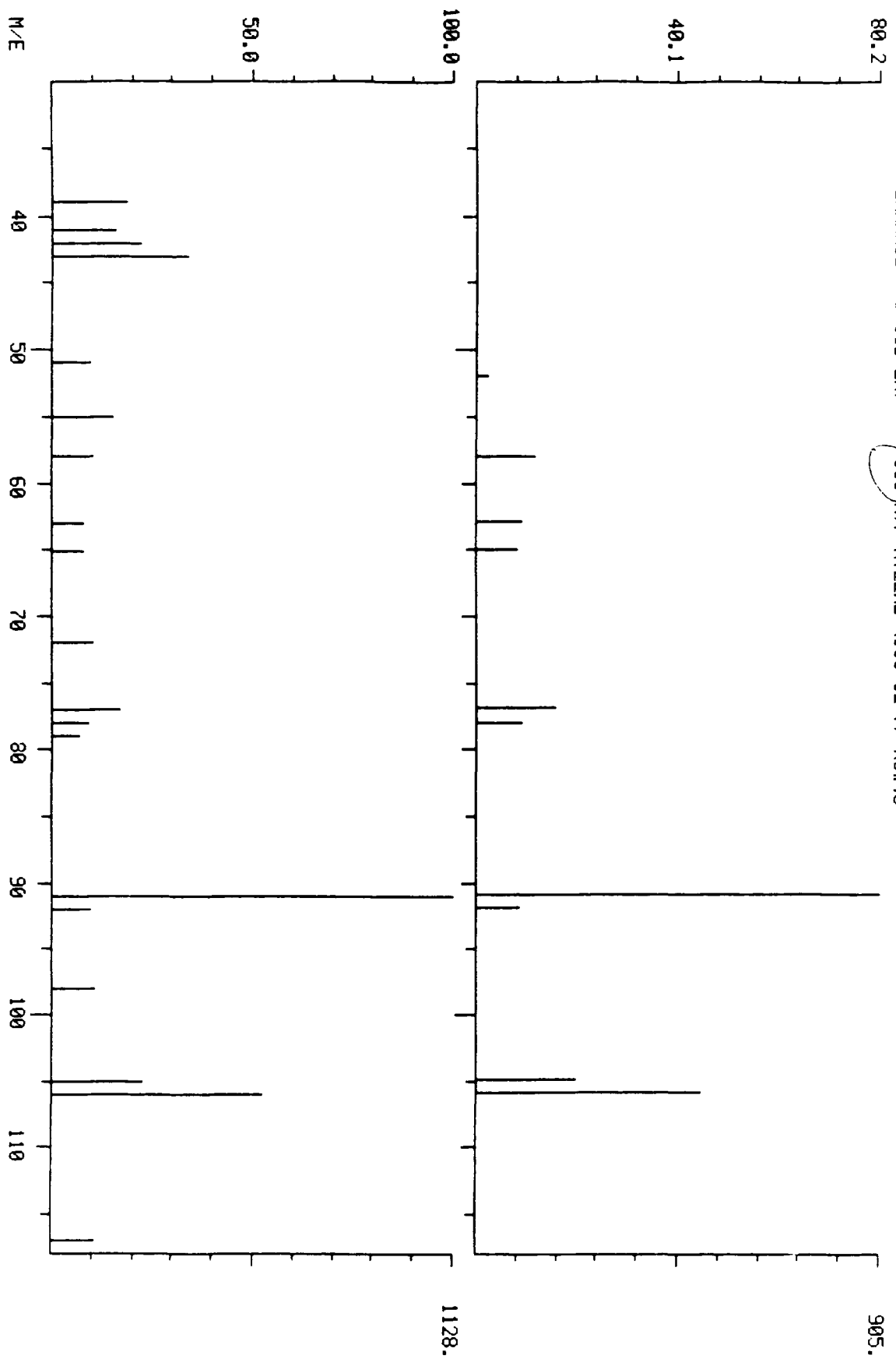


C8-H10
Y WT 11.05
B PK 91
RANK 1
IN 48
PUR 803

DUAL MASS SPECTRUM
12/22/89 6:56:00 + 13:04
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (5 158 2N) (330) M,P-XYLENE (133-02-7) R0#48

COMPUchem LABS

DATA: GH009588C19 #1046 BASE M/E: 91 / 91
RIC: 2363. / 4391.



LIBRARY SEARCH
12/22/89 6:56:00 + 13:00
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#8202A ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009688C19 #1040

BASE M/E: 42
RIC: 4375.

1595
SAMPLE

06.H10.0
1595
M WT 98
B PK 42
RANK 1
IN 895
PUR 848

CYCLOPENTANONE, 2-METHYL- CAS# 1128-72-5

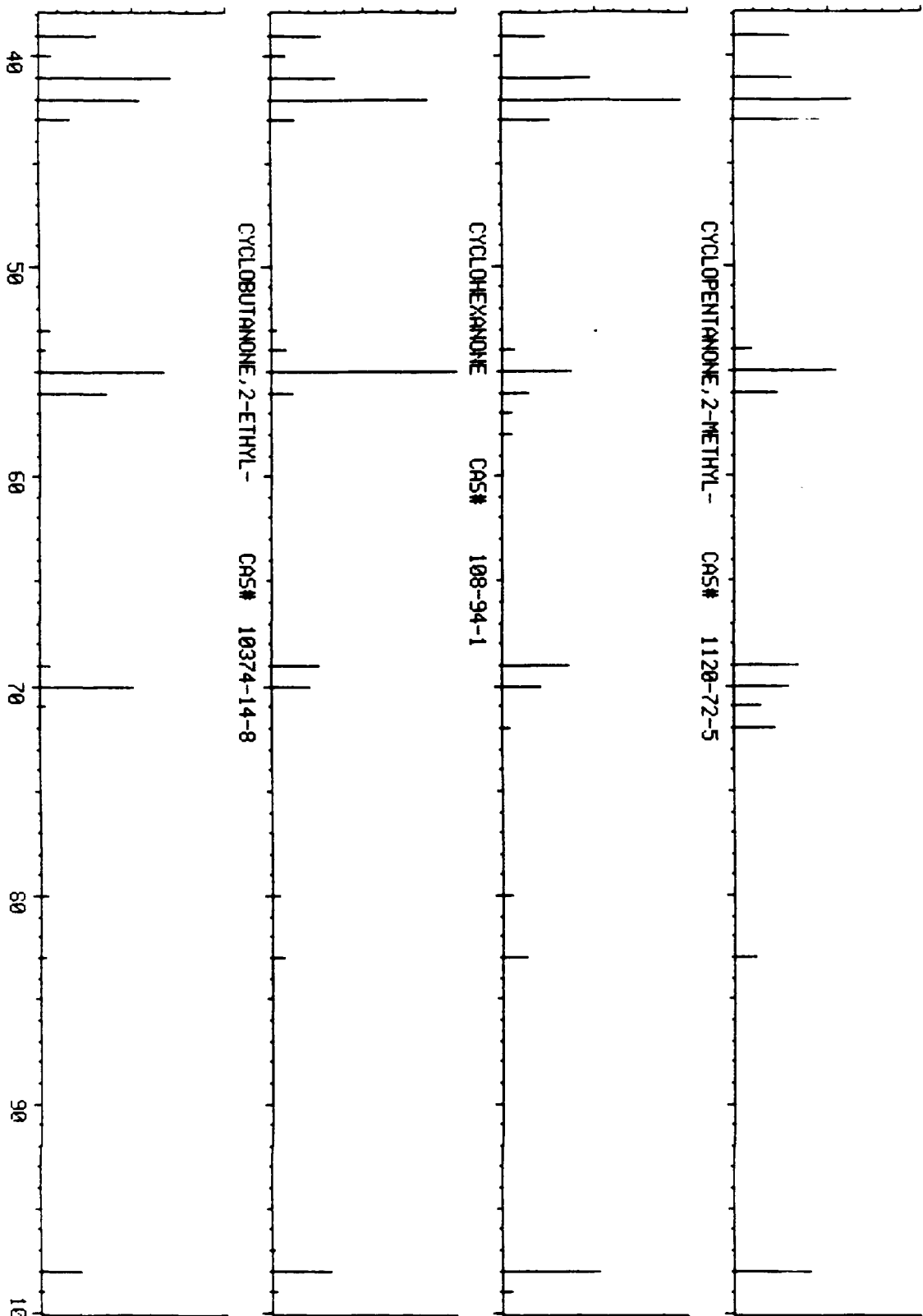
06.H10.0
1595
M WT 98
B PK 55
RANK 2
IN 876
PUR 795

CYCLOHEXANONE CAS# 108-94-1

06.H10.0
1595
M WT 98
B PK 41
RANK 3
IN 915
PUR 766

CYCLOBUTANONE, 2-ETHYL- CAS# 10374-14-8

M/E 40 50 60 70 80 90 100

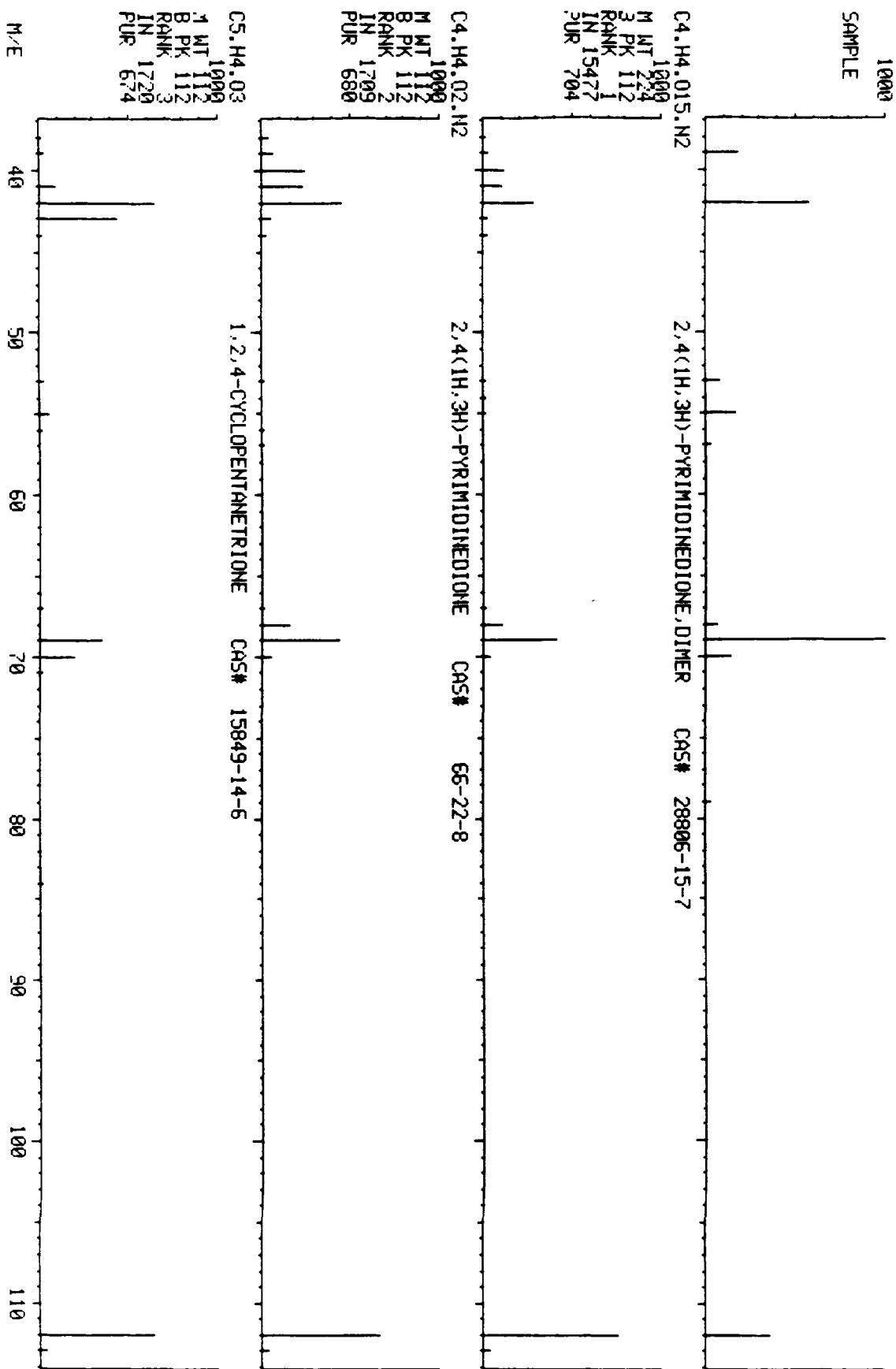


LIBRARY SEARCH
12/22/89 6:56:00 + 14:22
SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A ON#19
ENHANCED (5 156 2H 0T)

COMPUCHEM LABS

DATA: GH009688C19 #1150

BASE M/E: 69
RIC: 2387.



LIBRARY SEARCH
12/22/89 5:56:00 + 15:34
SAMPLE: 5G CC#309686 CASE#18756.7 EPA#B202A ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009688C19 #1246

BASE M/E: 84
RIC: 4343.

1549
SAMPLE

C7.H12.0

M WT 1549
B PK 112
RANK 84
IN 1787
PUR 741

CYCLOPENTANONE, 2-ETHYL-

CAS# 4971-18-0

C7.H12.0

M WT 1549
B PK 112
RANK 55
IN 1773
PUR 620

CYCLOHEXANONE, 4-METHYL-

CAS# 589-92-4

C8.H16

M WT 1549
B PK 112
RANK 53
IN 1854
PUR 620

1-PENTENE, 3-ETHYL-2-METHYL-

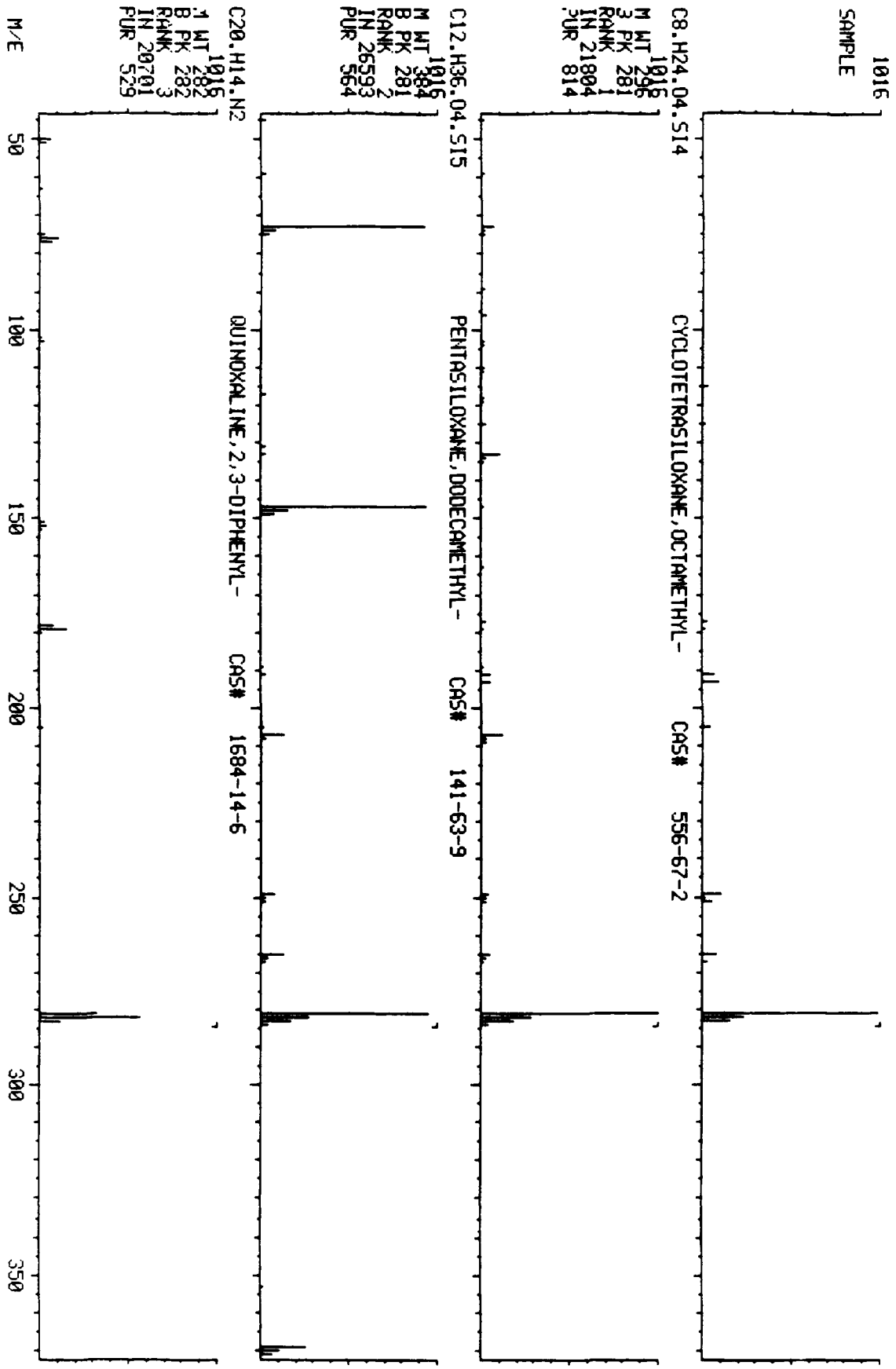
CAS# 19780-66-6



LIBRARY SEARCH
 12/22/89 6:56:00 + 15:55
 SAMPLE: 5G CC#309688 CASE#18756.7 EPA#B202A DN#19
 ENHANCED (5 158 2M 0T)

COMPUCHEM LABS
 DATA: GH009688C13 #1274

BASE M/E: 281
 RIC: 4863.



LAB INSTRUCTIONS:

SAMPLED DATE 12/18/89

RECEIPT DATE 12/20/89

CASE#: 18756 7

DUE DATE:

VOA
GC/MS WORKSHEET

COMPUCHEM#: 309688

R1 [] R2 [] D1 [] ()
R3 [] R4 [] D2 [] ()

L.L. SOLID, EPA SOW 2/88

Sample Prep Code---155
Instrument Code----413
Compound List-----494
Surrogate Std-----394
Internal Std-----036

=====

SAMPLE ID#:	B202A	Dry Wt. Factor	<u>1.14</u>	% Moisture	<u>12</u>
-------------	-------	----------------	-------------	------------	-----------

=====

GC/MS ANALYSIS

Amount Purged: [/] 10mls/Xg soil or [] Dilution _____ ul/100ul/Xg soil
Internal Standard Volume Added 5 ul
Surrogate Standard Volume Added 5 ul
BFB Filename 06891721019 Disk (2956A)
Blank Filename 614009785019 Disk ()
Standard Filename 65891777019 Disk ()
Sample Filename 64009698019 Disk ()

DEC 28 1989
GC/MS - UP 187-16

ANALYST(S): Injection 187-16

GC/MS REVIEW

CONDITION
CODE

OK

Entry Codes OK, EA, ES, SM, JS, SL, SH, JA, DA

Non-Entry Codes IM, IL, IH, SW, CT, CS, PC, NR
IF, LA, DI, CO, RN, DW, SI, SF
UP, BB, OT, VC, FO, NS

Extraneous Peak Search Results:

of Peaks Found: 4

Disposition: [/] Complete

[] Reprep neat required

[] Reprep using _____ g

Quality Assurance Notice(s):

Notices Required 1

[] Dilute ()

COMMENTS:

GC/MS Review OK Date 12/27/89 Auditor [Signature] Date 12/28/89

REPORT INTEGRATION

Final Reportable Package(s): GHO-C19

Total # of Injections: 1

QA COMMENTS:

Initials _____ Date ____/____/____

FINAL REVIEW

Initials _____ Date ____/____/____
AC1007 (05/89)

VOLATILE PREPARATION WORKSHEET

QUEUE # 19 PREP CODE -155 ASSIGNED TO Hushad Jeshi DATE 12/20/89

SAMPLE NUMBER	CASE NUMBER	SDG	QC SAMPLE		SAMPLE WEIGHT (G) VOLUME (ML)	SAMPLE ID	COMMENTS
			TYPE	ORIGINAL			
309763	18520	0164			5.0g	TB-12	
309768	↓	↓			5.0g	WB5-1	
309769	↓	↓			5.0g	WB5-2	
309770	↓	↓			5.0g	WB5-4	
309679	18756	0007			5.0g	B201A	
309680	↓	↓	SS	309679	5.0g	B201A MS	
309681	↓	↓	SS	309679	5.0g	B201A MSD	
309682	↓	↓	BS		0.0g		
309686	↓	↓			5.0g	B201B	
309687	↓	↓			5.0g	B202AR	
309688	↓	↓			5.0g	B202A	
309689	↓	↓			5.0g	B202B	
309690	↓	↓			5.0g	B202C	
309784			B1		5.0 ml	B3	
309785			B2		0.0 ml	B4	
309786			B3		0.0 ml	B5	
309787			B4		0.0 ml	B6	
309788			B5		0.0 ml	B7	

SURROGATE # 4 LOT # 1 MANUAL OPERATOR 739 / 516
 AMOUNT _____
 RELINQUISHED BY [Signature] DATE 12/20/89 RECEIVED BY [Signature] DATE 12/20/89

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	467	49300	50.0		X) 14
221	50	CHLOROMETHANE				BDL	10
231	62	VINYL CHLORIDE				BDL	10
220	94	BROMOMETHANE				BDL	10
209	64	CHLOROETHANE				BDL	10
216	96	1,1-DICHLOROETHENE				BDL	6
254	76	CARBON DISULFIDE				BDL	
252	43	ACETONE (2-PROPANONE)			32.6	33 32.6	10
248	114 I	1,4-DIFLUOROBENZENE (IS)	611	196000	50.0		
222	84	METHYLENE CHLORIDE			29.3	29 33.8	5
226	96	TRANS-1,2-DICHLOROETHENE				BDL	5
214	63	1,1-DICHLOROETHANE				BDL	5
257	43	VINYL ACETATE				BDL	10
237	96	CIS-1,2-DICHLOROETHENE				BDL	5
253	72	2-BUTANONE			22.9	23 26	10
211	83	CHLOROFORM				BDL	5
227	97	1,1,1-TRICHLOROETHANE				BDL	5
206	117	CARBON TETRACHLORIDE				BDL	5
203	78	BENZENE				BDL	5
215	62	1,2-DICHLOROETHANE				BDL	5
270	117 I	D5-CHLOROBENZENE (IS) RO#29	999	206000	50.0		
229	130	TRICHLOROETHENE				BDL	5
217	63	1,2-DICHLOROPROPANE				BDL	5
212	83	BROMODICHLOROMETHANE				BDL	5
218	75	CIS-1,3-DICHLOROPROPENE				BDL	5
256	43	4-METHYL-2-PENTANONE			2.2	BDL 2.2	10
225	92	TOLUENE			1.3	1.3 1.7	5
250	75	TRANS-1,3-DICHLOROPROPENE				BDL	5
228	97	1,1,2-TRICHLOROETHANE				BDL	5
224	164	TETRACHLOROETHENE				BDL	5
255	43	2-HEXANONE			11.1	11 13	10
208	129	DIBROMOCHLOROMETHANE				BDL	5
207	112	CHLOROBENZENE				BDL	5
219	106	ETHYLBENZENE			2.4	BDL 2.4	5
330	106	M, P-XYLENE			1.7	1.7 2.2	5
239	106	O-XYLENE				BDL	5
251	104	STYRENE				BDL	5
205	173	BROMOFORM				BDL	5
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	5
259	65 S	D4-1,2-DICHLOROETHANE RO#57			48.4	97.%	
247	95 S	BROMOFLUOROBENZENE			46.1	92.%	
233	98 S	D8-TOLUENE RO#59			47.6	95.%	
289	106	XYLENES (TOTAL)			1.7	2.2 2.7	5

CORRECTED/REVIEWED BY *C. J. ...*
(GC/MS DATA REVIEWER)

DATE 12/27/89

CMP	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
299	96	1,2-DICHLOROETHENE (TOTAL)				BDL	
CHECKSUMS:							
		3979.	2077	451300.		397.4	389.

CORRECTED/REVIEWED BY *C. J. St...*
(GC/MS DATA REVIEWER)

DATE 12-27-89

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P
40	258	D4-1,2-DICHLOROETHANE RO#57	48.4	50.0	97.	70-121	X
41	247	BROMOFLUOROBENZENE	46.1	50.0	92.	74-121	X
42	233	D8-TOLUENE RO#59	47.6	50.0	95.	81-117	X

* ADVISORY SURROGATE ONLY

++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

CORRECTION FACTOR CALCULATION:

$$\frac{5.0 \text{ G}}{\text{WET WEIGHT OF SAMPLE (G)}} \times \frac{\text{GC/MS DILUTION FACTOR}}{\text{GC/MS DILUTION FACTOR}} \times \frac{\text{DRY WEIGHT FACTOR}}{\text{DRY WEIGHT FACTOR}} =$$

$$\frac{5.0 \text{ G}}{5.00 \text{ (G)}} \times \frac{1.0}{1.0} \times \frac{1.14}{1.00} = 1.14$$

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:

THE SURROGATES ARE ADDED TO THE SAMPLE PRIOR TO SPARGING.
SURROGATE SPIKE CONVERSION FACTOR = 1.

VERSION 8

CORRECTED/REVIEWED BY EK Stebbins
(GC/MS DATA REVIEWER)

DATE 12-27-89

LABORATORY NOTICE

Sample I.D. # 309688

Client I.D. # B202A

CASE # 18756.7

The analysis of the volatile fraction of this sample indicated the presence of one or more extraneous peaks at the scan(s) listed below which were tentatively identified as siloxanes. This is a frequently observed laboratory artifact and should not be considered an actual sample constituent. It is being reported as an instrument artifact. While the peak(s) have been included in the total number of Library Searches, they have not been counted as one of the ten most intense TIC's requiring Library Search. If the siloxane was also detected in the associated blank, it has been flagged in the sample data with a "B", whether or not its concentration in the blank was great enough to require a Library Search. The data is being reported with reference to this qualifier.

scan(s): 1274 _____

Data Reviewer: OK [Signature]

Date: 12 27 88

Approved by: Robert J. Whitehead
Manager, Quality Assurance

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009689C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 14 Date Analyzed: 12/22/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	12	U
74-83-9	-----Bromomethane	12	U
75-01-4	-----Vinyl Chloride	12	U
75-00-3	-----Chloroethane	12	U
75-09-2	-----Methylene Chloride	30	B
67-64-1	-----Acetone	53	B
75-15-0	-----Carbon Disulfide	6	U
75-35-4	-----1,1-Dichloroethene	6	U
75-34-3	-----1,1-Dichloroethane	6	U
540-59-0	-----1,2-Dichloroethene (total)	6	U
67-66-3	-----Chloroform	6	U
107-06-2	-----1,2-Dichloroethane	6	U
78-93-3	-----2-Butanone	40	U
71-55-6	-----1,1,1-Trichloroethane	6	U
56-23-5	-----Carbon Tetrachloride	6	U
108-05-4	-----Vinyl Acetate	12	U
75-27-4	-----Bromodichloromethane	6	U
78-87-5	-----1,2-Dichloropropane	6	U
10061-01-5	-----cis-1,3-Dichloropropene	6	U
79-01-6	-----Trichloroethene	6	U
124-48-1	-----Dibromochloromethane	6	U
79-00-5	-----1,1,2-Trichloroethane	6	U
71-43-2	-----Benzene	6	U
10061-02-6	-----Trans-1,3-Dichloropropene	6	U
75-25-2	-----Bromoform	6	U
108-10-1	-----4-Methyl-2-Pentanone	12	U
591-78-6	-----2-Hexanone	23	U
127-18-4	-----Tetrachloroethene	6	U
79-34-5	-----1,1,2,2-Tetrachloroethane	6	U
108-88-3	-----Toluene	15	U
108-90-7	-----Chlorobenzene	6	U
100-41-4	-----Ethylbenzene	12	U
100-42-5	-----Styrene	3	J
1330-20-7	-----Total Xylenes	54	J

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202B

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309689
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009689C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 14 Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 7

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

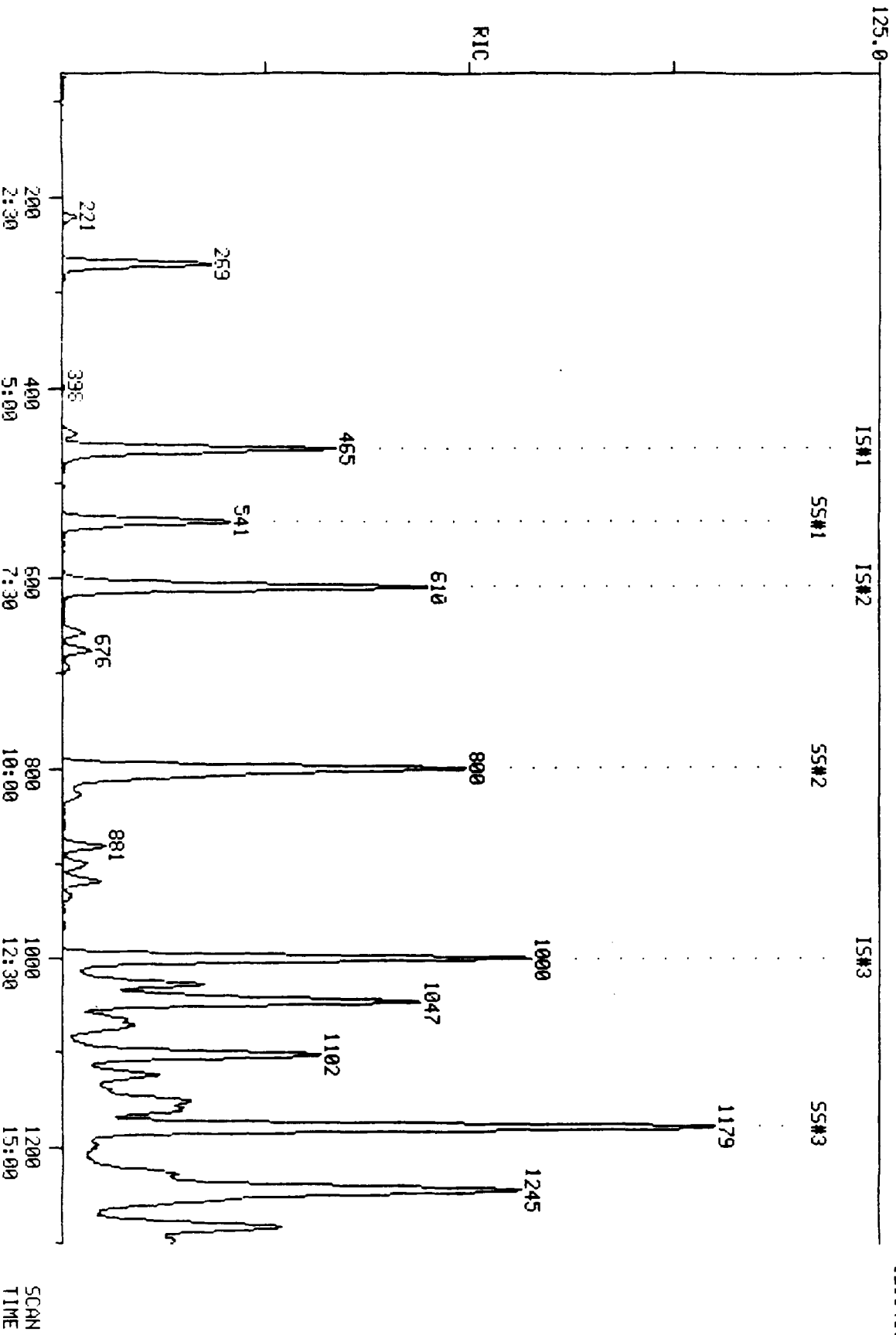
CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	13.39	8.1	J
2.	UNKNOWN	14.05	7.0	J
3.	UNKNOWN	14.40	8.1	J
4.	UNKNOWN	14.49	8.1	J
5. 103-65-1	BENZENE, PROPYL-	15.35	8.1	J
6. 620-14-4	BENZENE, 1-ETHYL-3-METHYL-	15.55	84	J
7. 611-14-3	BENZENE, 1-ETHYL-2-METHYL-	16.04	21	J

RIC
12/22/89 7:40:00
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B DN#19
COND.:

COMPUCHEM LABS

COMPUCHEM DATA: GH009689C19 SCANS 71 TO 1300

123040.



QUANTITATION REPORT FILE: GH009689C19
 DATA: GH009689C19.TI
 12/22/89 7:40:00
 SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B ON#19
 CONDS.:
 SUBMITTED BY: 19 ANALYST: 1422

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
 RESP. FAC. FROM LIBRARY ENTRY -

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1
2	221 CHLOROMETHANE <74-87-3> RO#2
3	231 VINYL CHLORIDE <75-01-4> RO#3
4	220 BROMOMETHANE <78-83-9> RO#4
5	209 CHLOROETHANE <75-00-3> RO#5
6	216 1,1-DICHLOROETHENE <75-35-4> RO#8
7	254 CARBON DISULFIDE <75-15-0> RO#9
8	252 ACETONE (2-PROPANONE) <67-64-1> RO#13
9	*249 1,4-DIFLUOROBENZENE (IS) <540-36-3> RO#14
10	222 METHYLENE CHLORIDE <75-09-2> RO#16
11	226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17
12	214 1,1-DICHLOROETHANE <75-34-3> RO#19
13	257 VINYL ACETATE <108-05-4> RO#20
14	237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21
15	253 2-BUTANONE <78-93-3> RO#22
16	211 CHLOROFORM <67-66-2> RO#23
17	227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24
18	206 CARBON TETRACHLORIDE <56-23-5> RO#25
19	203 BENZENE <71-43-2> RO#26
20	215 1,2-DICHLOROETHANE <107-06-2> RO#27
21	*270 D5-CHLOROENZENE (IS) RO#29
22	229 TRICHLOROETHENE <79-01-6> RO#30
23	217 1,2-DICHLOROPROPANE <78-87-5> RO#31
24	212 BROMODICHLOROMETHANE <75-27-4> RO#33
25	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35
26	256 4-METHYL-2-PENTANONE <108-01-1> RO#36
27	225 TOLUENE <108-88-3> RO#37
28	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38
29	228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39
30	224 TETRACHLOROETHENE <127-18-4> RO#41
31	255 2-HEXANONE <591-78-6> RO#42
32	208 DIBROMOCHLOROMETHANE <124-48-1> RO#43
33	207 CHLOROBENZENE <108-90-7> RO#45
34	219 ETHYLBENZENE <100-41-4> RO#47
35	330 M,P-XYLENE <133-02-7> RO#48
36	239 O-XYLENE <133-02-7> RO#49
37	251 STYRENE <100-42-5> RO#50
38	205 BROMOFORM <75-25-2> RO#51
39	223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54
40	*258 D4-1,2-DICHLOROETHANE RO#57
41	*247 BROMOFLUOROBENZENE <460-00-4> RO#58
42	*233 D8-TOLUENE RO#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	465	5:49	1	1.000	A BB	46782.	50.000 UG/KG	10.15
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	221	2:46	1	0.475	A BB	11289.	45.597 UG/KG	9.26
9	114	609	7:37	9	1.000	A BB	191540.	50.000 UG/KG	10.15
10	84	270	3:22	1	0.581	A BB	40013.	26.024 UG/KG	5.28
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	448	5:36	1	0.963	A BB	3418.	34.087 UG/KG	6.92
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	NOT FOUND							
20	62	NOT FOUND							
21	117	1000	12:30	21	1.000	A BB	203559.	50.000 UG/KG	10.15
22	130	NOT FOUND							
23	63	NOT FOUND							
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	799	9:59	21	0.799	A BB	2497.	2.390 UG/KG	0.49
27	92	808	10:06	21	0.808	A BB	38527.	12.821 UG/KG	2.60
28	75	NOT FOUND							
29	97	NOT FOUND							
30	164	NOT FOUND							
31	43	919	11:29	21	0.919	A VB	16058.	20.205 UG/KG	4.10
32	129	NOT FOUND							
33	112	NOT FOUND							
34	106	1027	12:50	21	1.027	A BV	20927.	10.143 UG/KG	2.06
35	106	1047	13:05	21	1.047	A VB	73966.	25.787 UG/KG	5.23
36	106	1103	13:47	21	1.103	A BB	53160.	20.995 UG/KG	4.26
37	104	1106	13:49	21	1.106	A BB	12112.	2.707 UG/KG	0.55
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	540	6:45	1	1.161	A BB	94248.	48.275 UG/KG	9.80
41	95	1179	14:44	21	1.179	A BB	166628.	47.845 UG/KG	9.71
42	98	800	10:00	21	0.800	A BB	195344.	45.798 UG/KG	9.30

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:54	0.99	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:06		10.000			50.00		0.952	
3	1:11		10.000			50.00		1.256	
4	1:26		10.000			50.00		1.579	
5	1:34		10.000			50.00		0.796	
6	2:37		5.000			50.00		1.346	
7	2:46		5.000			50.00		3.709	
8	2:48	0.99	10.000	0.05	45.60	50.00	0.241	0.265	0.91
9	7:42	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	3:27	0.98	5.000	0.12	26.02	50.00	0.855	1.643	0.52
11	3:53		5.000			50.00		1.486	
12	4:34		5.000			50.00		2.635	
13	4:52		10.000			50.00		0.458	
14	5:33		5.000			50.00		1.675	

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:42	0.98	10.000	0.10	34.09	50.00	0.073	0.107	0.68
16	6:10		5.000			50.00		3.433	
17	6:18		5.000			50.00		0.814	
18	6:34		5.000			50.00		0.755	
19	6:53		5.000			50.00		0.895	
20	6:58		5.000			50.00		2.523	
21	12:34	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	8:01		5.000			50.00		0.569	
23	8:22		5.000			50.00		0.352	
24	8:55		5.000			50.00		0.701	
25	9:41		5.000			50.00		0.690	
26	10:04	0.99	15.000	0.05	2.39	50.00	0.012	0.257	0.05
27	10:10	0.99	5.000	0.16	12.82	50.00	0.189	0.738	0.26
28	10:42		5.000			50.00		0.342	
29	10:58		5.000			50.00		0.349	
30	11:06		5.000			50.00		0.723	
31	11:33	0.99	15.000	0.06	20.21	50.00	0.079	0.195	0.40
32	11:35		5.000			50.00		0.690	
33	12:37		5.000			50.00		1.092	
34	12:55	0.99	5.000	0.21	10.14	50.00	0.103	0.507	0.20
35	13:10	0.99	5.000	0.21	25.79	50.00	0.363	0.705	0.52
36	13:52	0.99	5.000	0.22	20.99	50.00	0.261	0.622	0.42
37	13:54	0.99	5.000	0.22	2.71	50.00	0.060	1.099	0.05
38	14:07		5.000			50.00		0.664	
39	15:16		5.000			50.00		0.557	
40	6:52	0.98	5.000	0.23	48.27	50.00	2.015	2.087	0.97
41	14:49	0.99	5.000	0.24	47.84	50.00	0.819	0.855	0.96
42	10:04	0.99	5.000	0.16	45.80	50.00	0.960	1.048	0.92

LIBRARY SEARCH
12/22/89 7:40:00 + 2:45
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009689C19 # 221

BASE M/E: 43
RIC: 2003.

1057
SAMPLE

C3.H6.O
M.WT 58
B.PK 43
RANK 1
IN 13
PUR 837

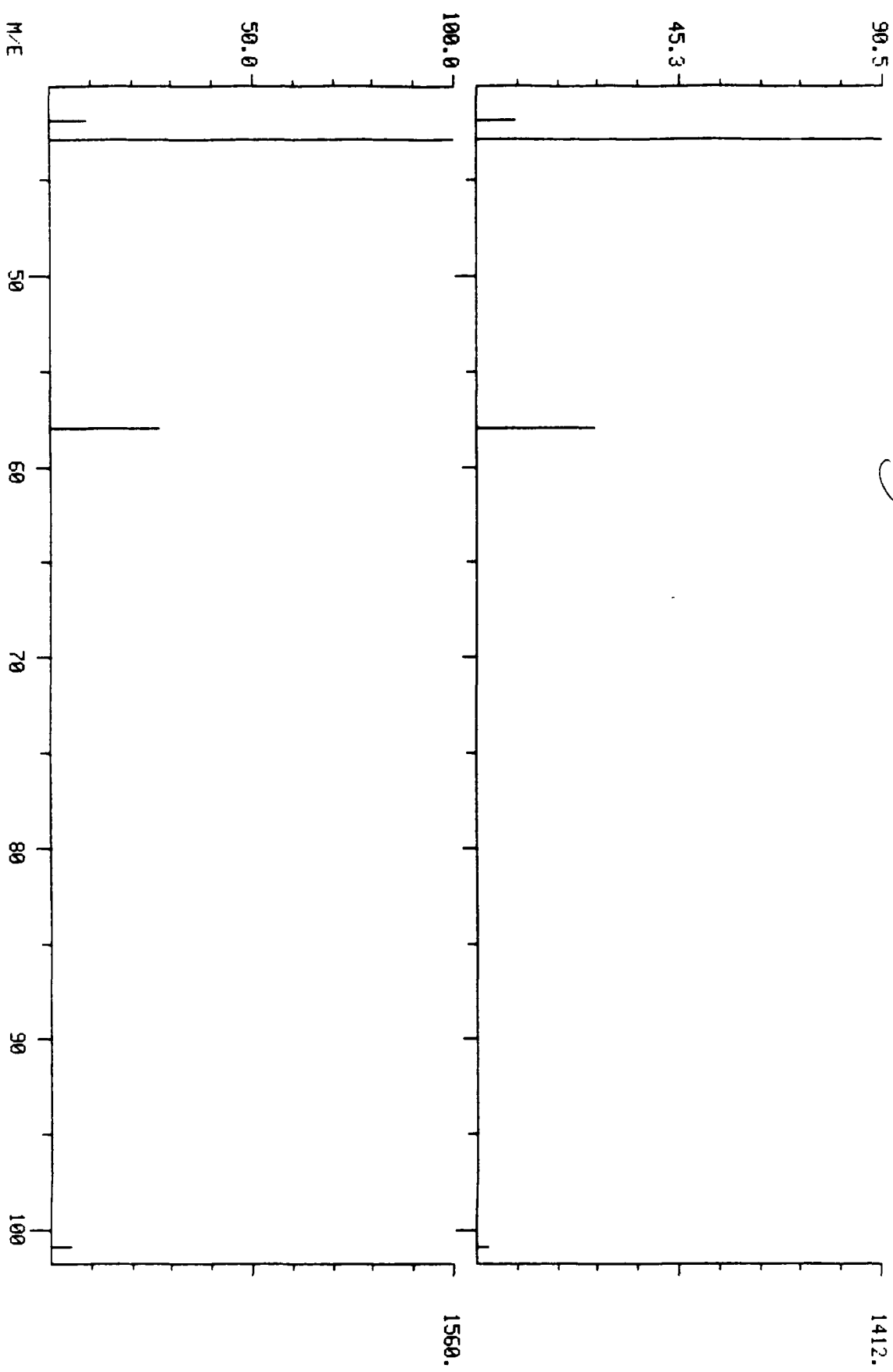
252 ACETONE (2-PROPANONE) <67-64-1> R0#13

1057

SAMPLE MINUS LIBRARY



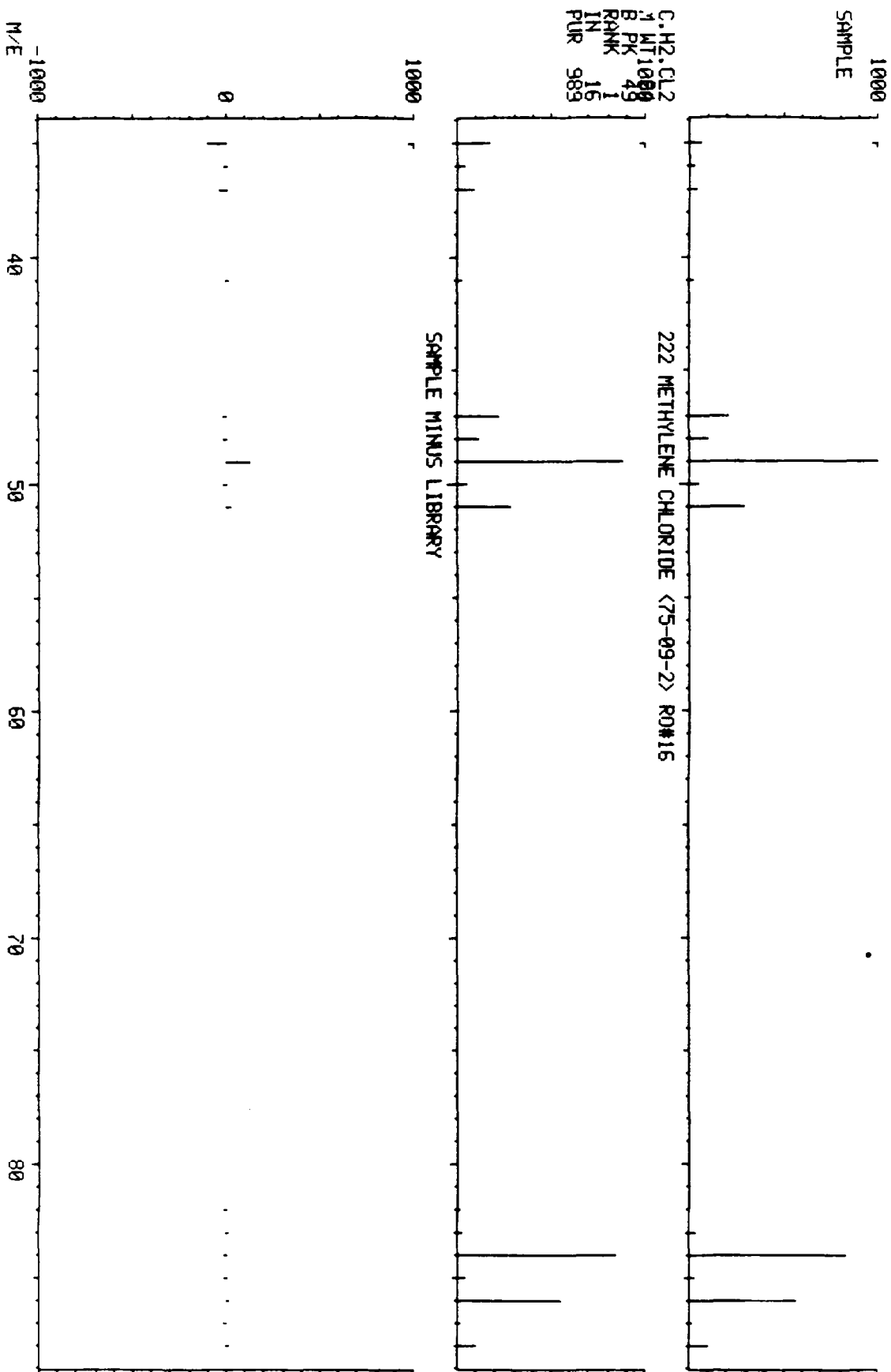
DUAL MASS SPECTRUM
12/22/89 7:40:00 + 2:46
SAMPLE: 5G CC#309689 CASE#18756.7 EPANR202B ON#19
ENHANCED (S 158 2N) (252) ACETONE (2-PROPANONE) <67-64-1> R0#13
COMPUCHEM LABS
DATA: GH009689C19 #221 BASE M/E: 43/ 43
RIC: 2003./ 2195.



LIBRARY SEARCH
12/22/89 7:40:00 + 3:22
SAMPLE: 5G.CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS
DATA: GH009689C19 # 270
BASE M/E: 49
RID: 21151.

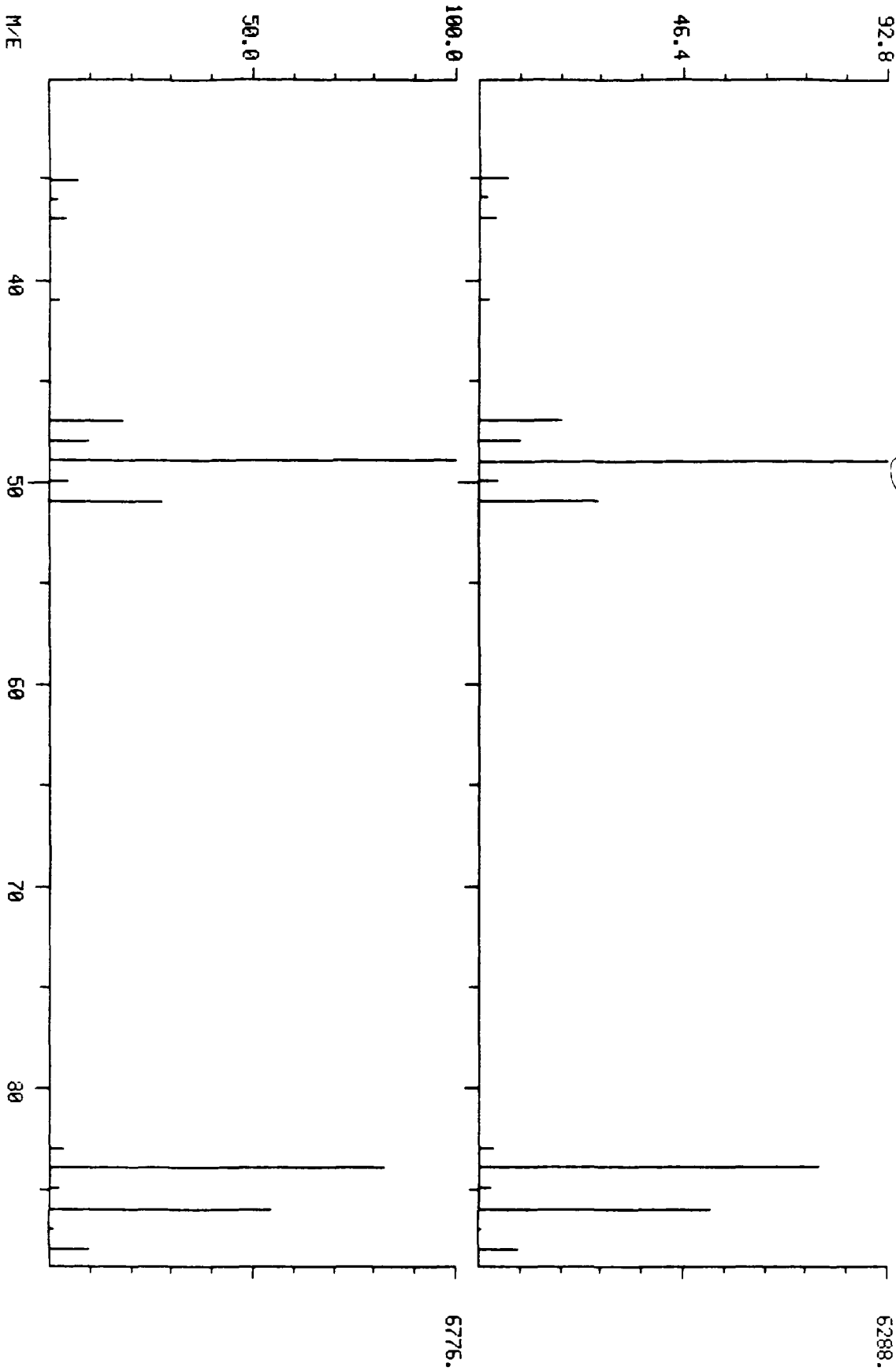
C.H2.CL2
1 MT 1000
8 PK 49
RANK 1
IN 16
PUR 989



DUAL MASS SPECTRUM
12/22/89 7:40:00 + 3:22
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B ON#19
ENHANCED (5 15B 2N) (222) METHYLENE CHLORIDE <75-09-2> RD#16

COMPUCHEM LABS

DATA: CH009689C19 #270 BASE M/E: 49/ 45
RIC: 21151.7 22143.





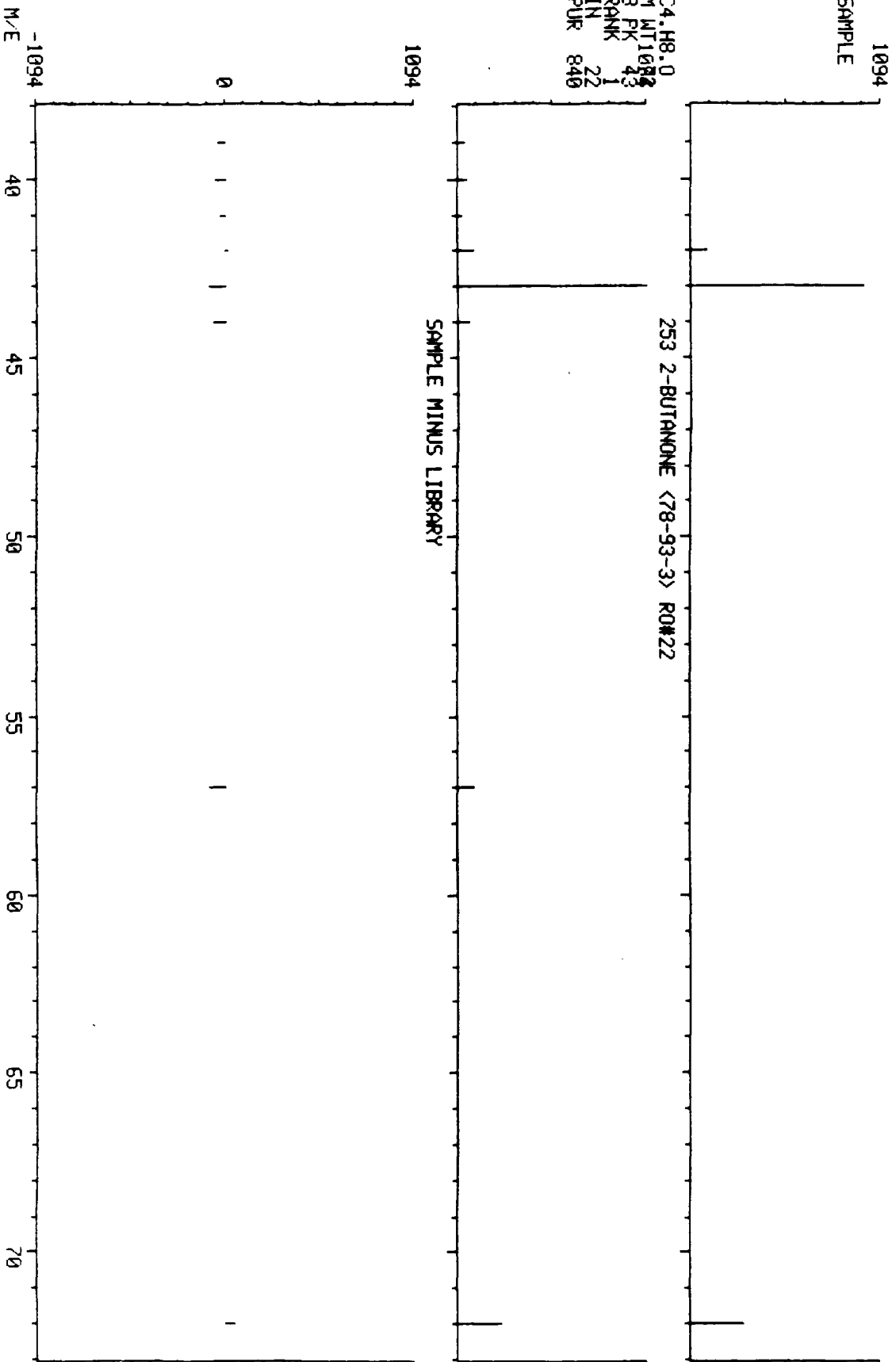
COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 7:40:00 + 5:36
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#R2028 ON#19
ENHANCED (5 15B 2N 0T)

COMPUCHEM LABS

DATA: GH009689C19 # 448

BASE M/E: 43
RIC: 1631.

C4.H8.O
1 MIT 1082
R PK 43
RANK 1
IN 22
PUR 840

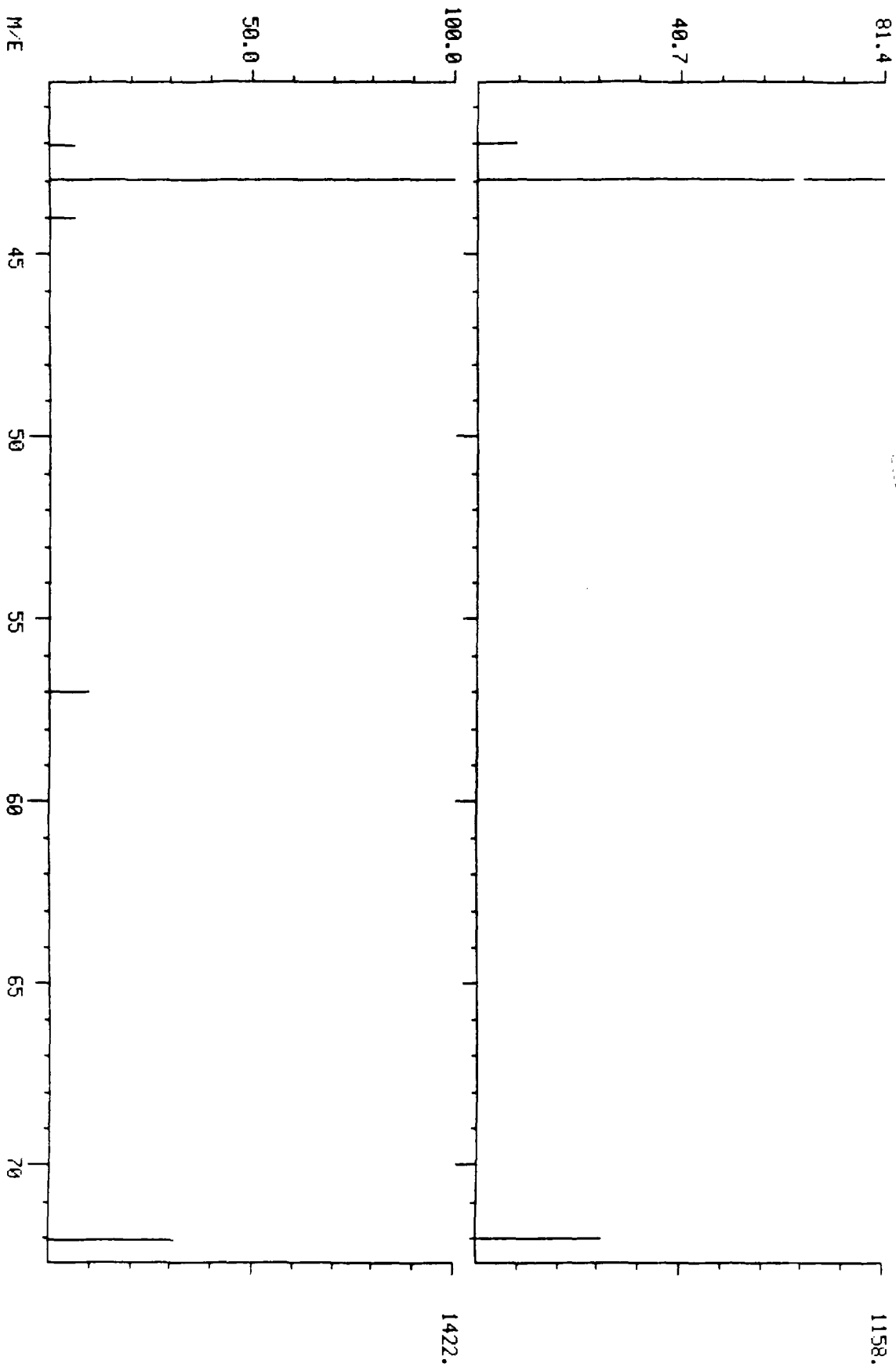


DUAL MASS SPECTRUM
12/22/89 7:40:00 + 5:36
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B ON#19
ENHANCED (5 156 2N) 253)2-BUTANONE (78-93-3) R0#22

COMPUCHEM LABS

DATA: CH009689C19 #448

BASE M/E: 43/ 43
RIC: 1631./ 2171.



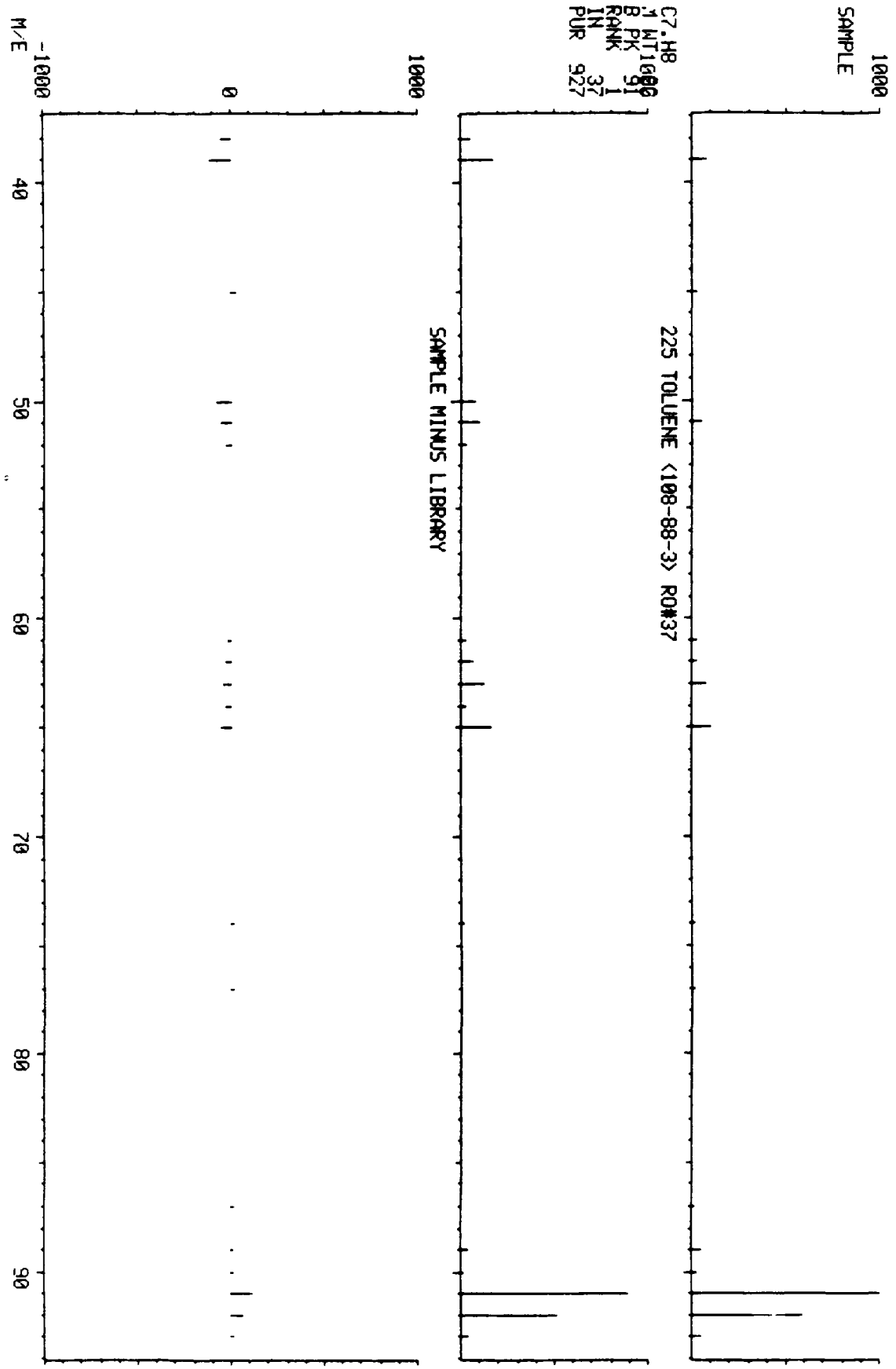


LIBRARY SEARCH
12/22/89 7:40:00 + 10:05
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (S 158 ZN 01)

COMPUchem LAB5
DATA: GH009689C19 # 808

BASE M/E: 91
RIC: 12559.

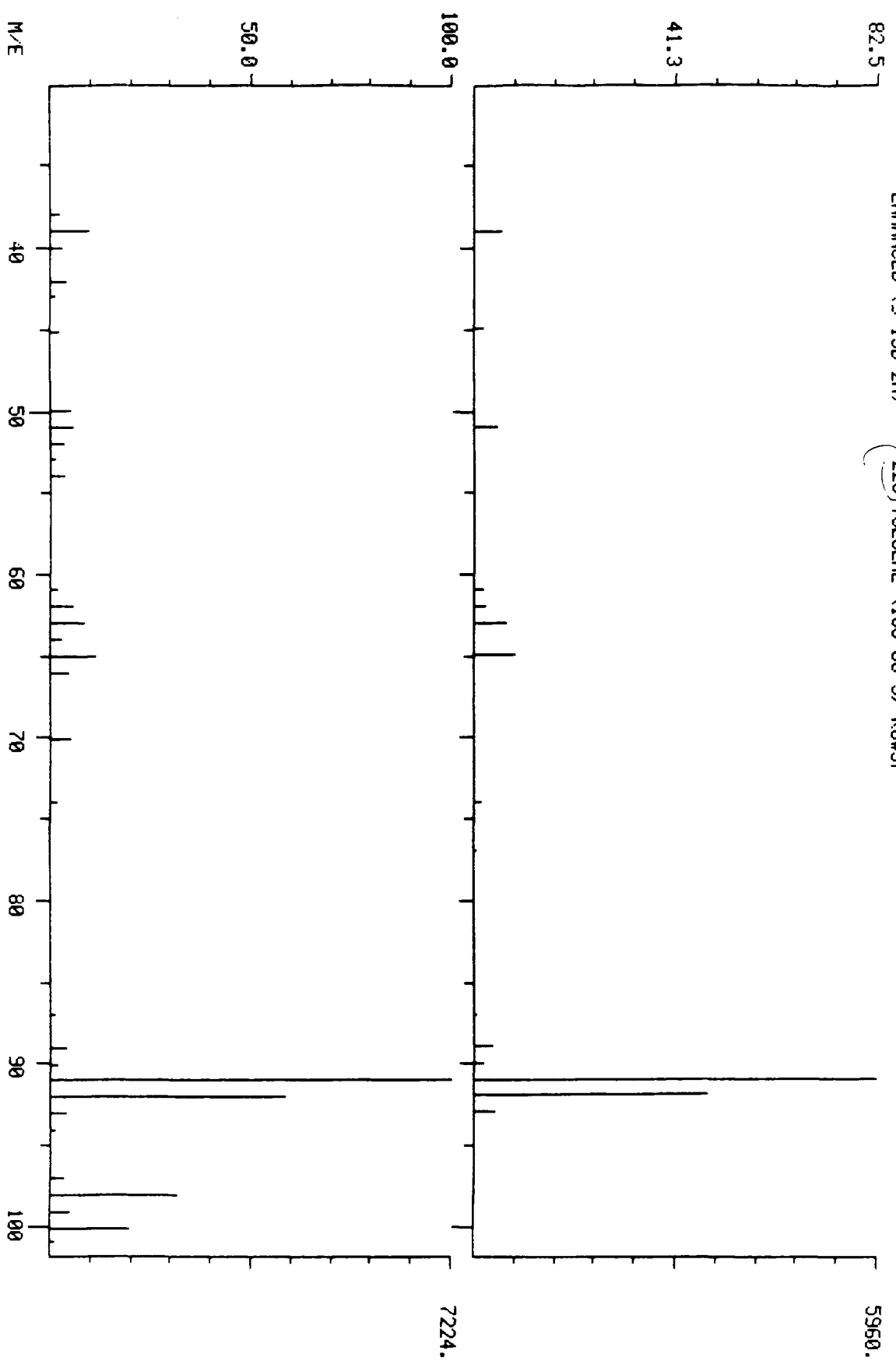
C7.H8
1 WT 1000
B PK 91
RANK 37
IN 1
PUR 927



DUAL MASS SPECTRUM
12/22/89 7:40:00 + 10:06
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B OH#19
ENHANCED (5 158 2M) (225) TOLUENE <108-88-3> RM#37

COMPUCHEM LABS

DATA: GH009689C19 #808 BASE M/E: 91 / 91
RIC: 12559. / 22591.

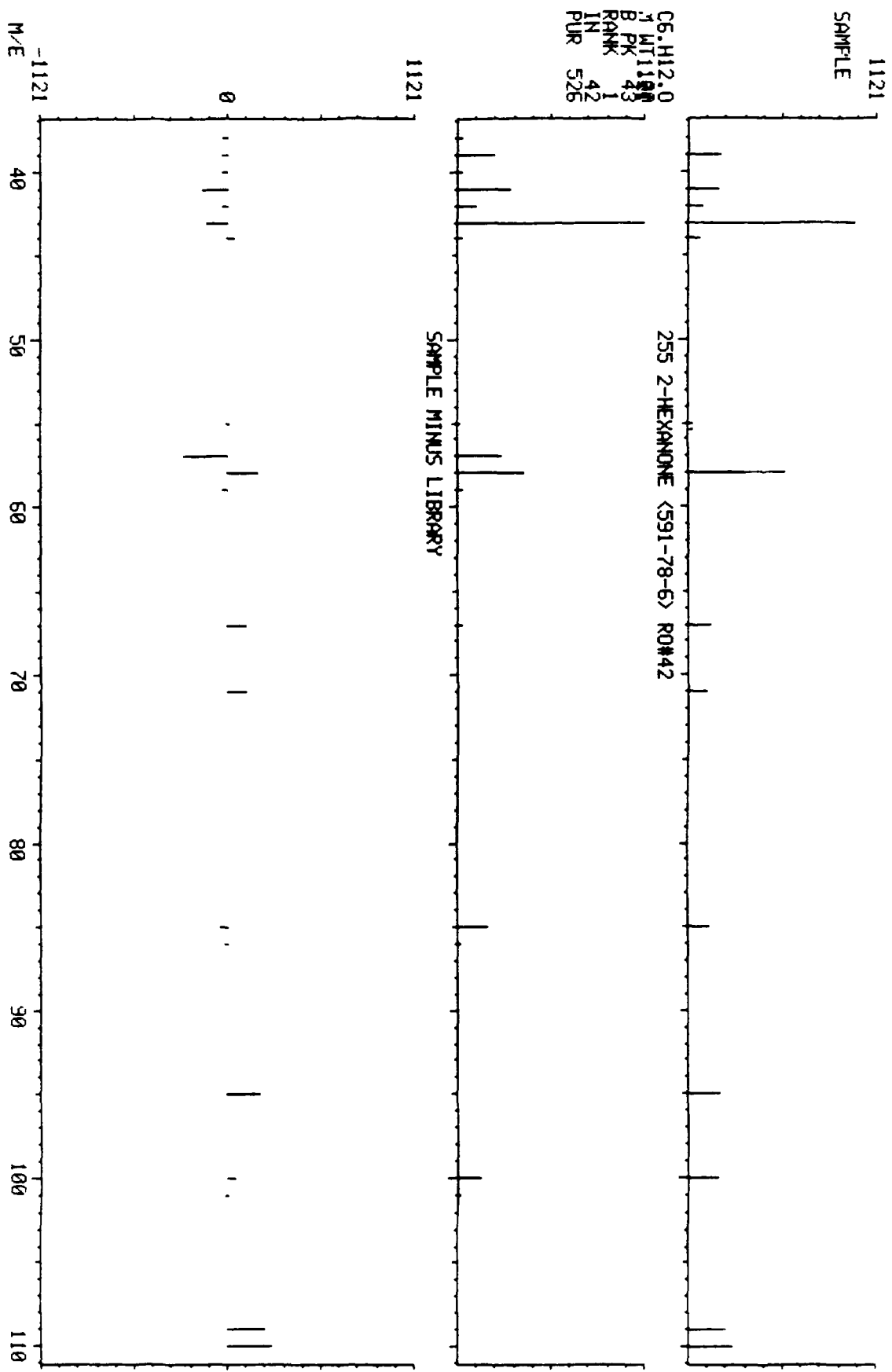


COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 7:40:00 + 11:29
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B ON#19
ENHANCED (5 158 2N 0T)

DATA: GH009689C19 # 919

BASE M/E: 43
RIC: 4359.

CG: 112.0
I: 111.0
B: 43
RANK: 1
IN: 42
PUR: 526

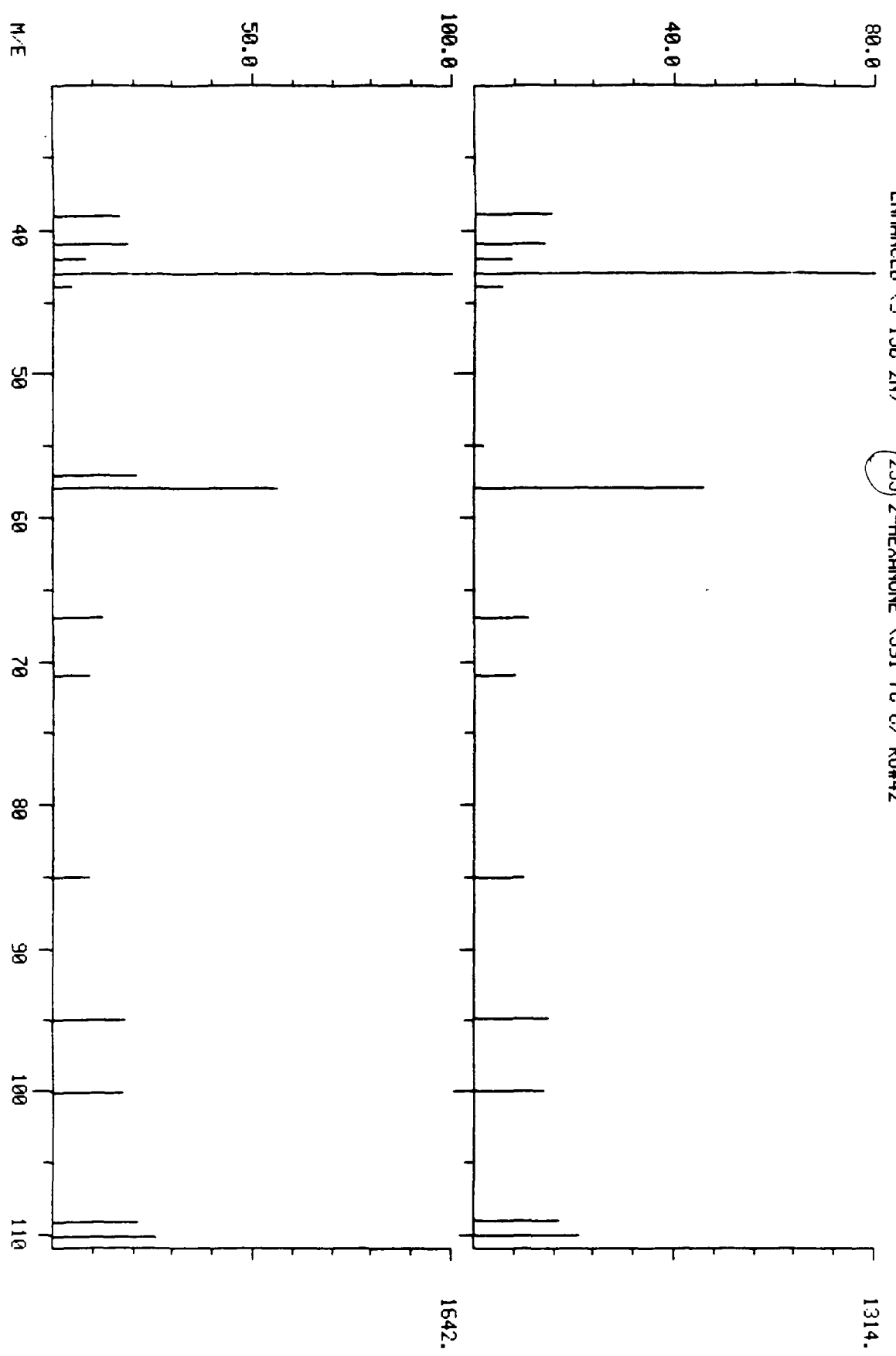


DUAL MASS SPECTRUM
12/22/89 7:40:00 + 11:29
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (5 15B 2N) (25) 2-HEXANONE (591-78-E) R0#42

COMPUCHEM LABS

DATA: GH009689C19 #919

BASE M/E: 43/ 43
RIC: 4359./ 5535.

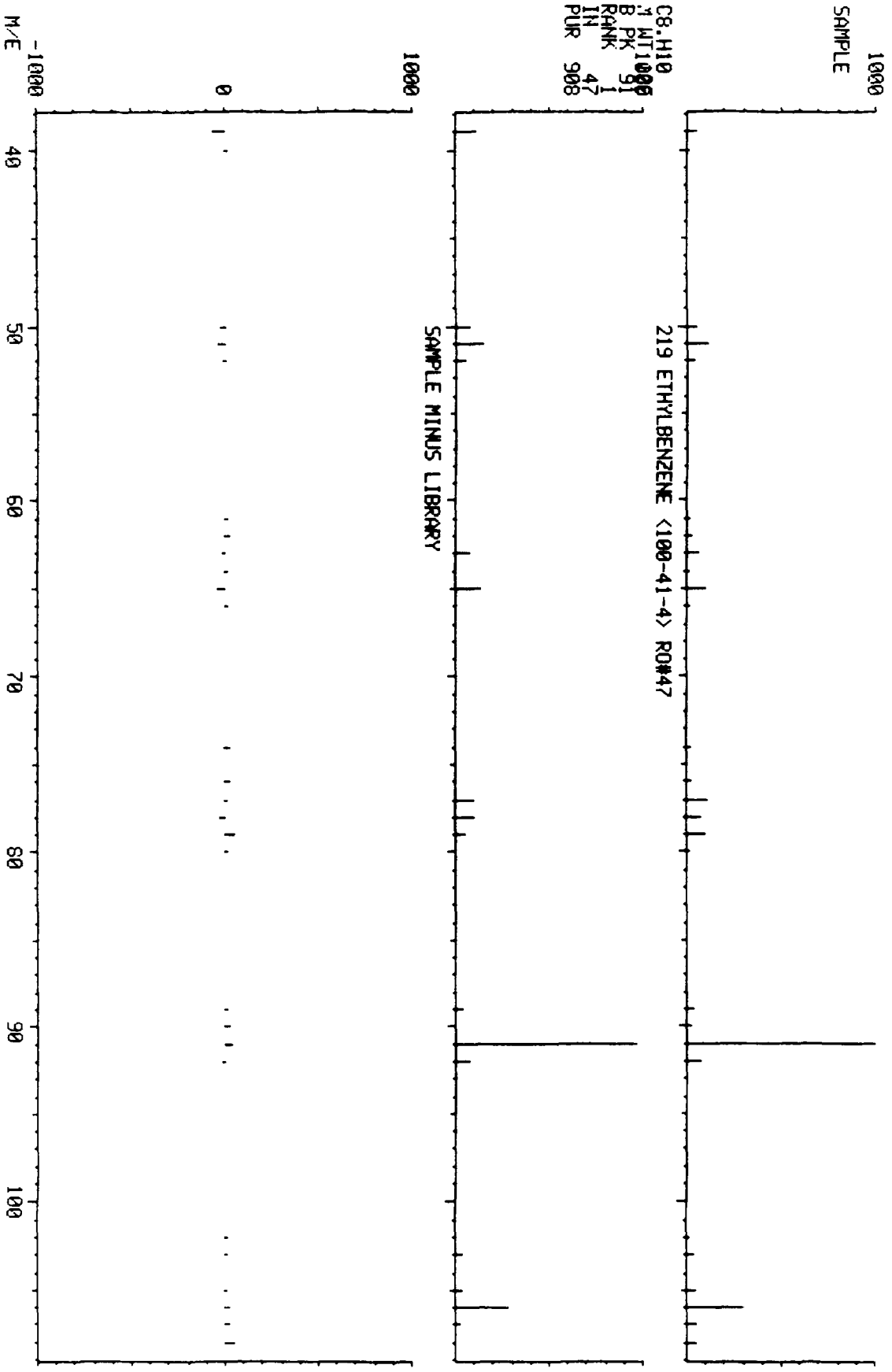


COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 7:40:00 + 12:50
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (5 15B 2N 0T)

DATA: CH009689C19 #1027

BASE M/E: 91
R1C: 15471.

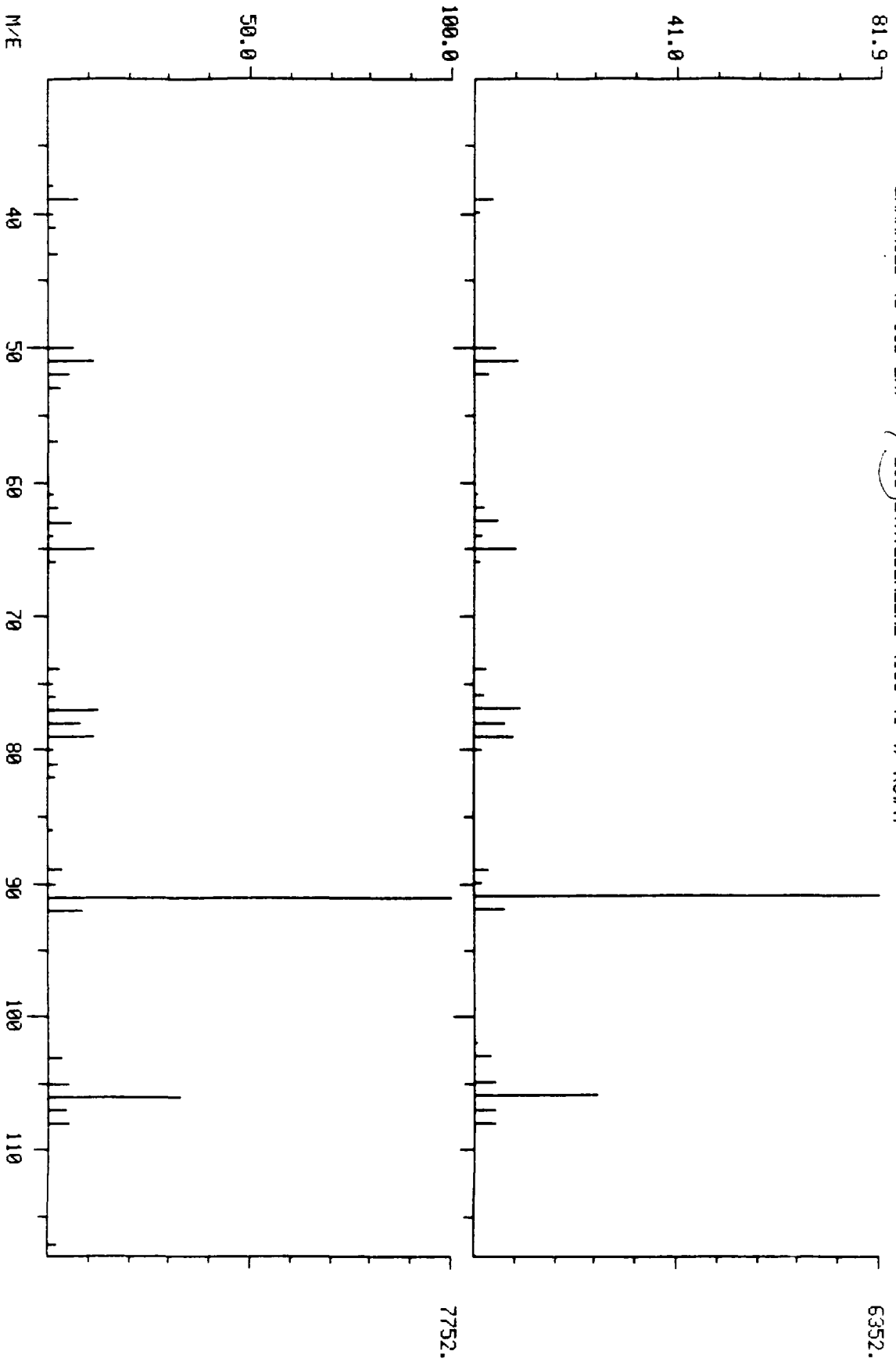
C8.H10
7 MT 1006
8 PK 91
RANK 47
IN 47
PUR 908



DUAL MASS SPECTRUM
12/22/89 7:40:00 + 12:50
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B ON#19
ENHANCED (5 15B 2N) 219 ETHYL BENZENE <100-41-4> R0#47

COMPUCHEM LABS

DATA: GH009689C19 #1027 BASE M/E: 91 / 91
RIC: 15471. / 21151.



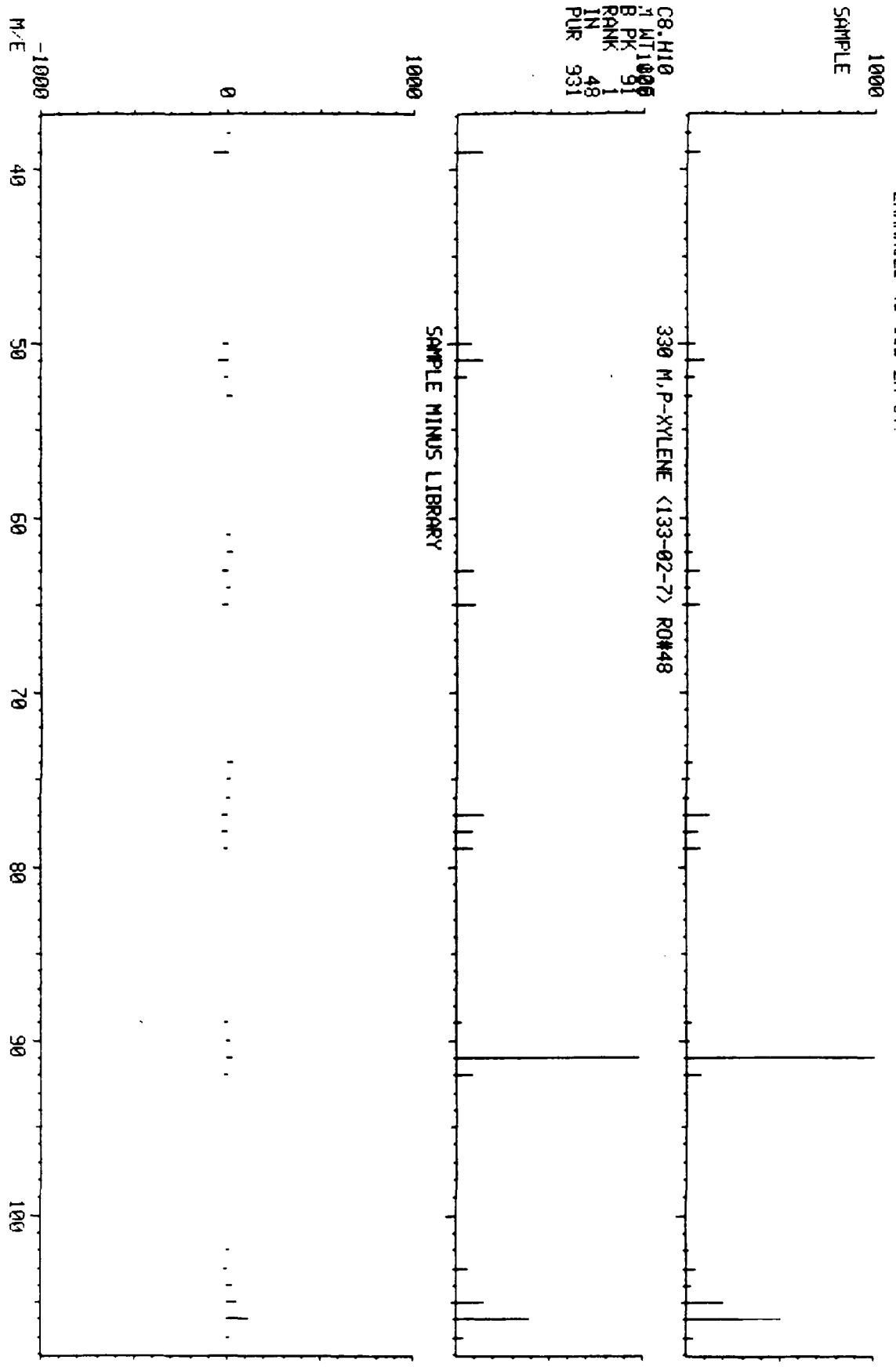
LIBRARY SEARCH
12/22/89 7:40:00 + 13:05
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

DATA: GH009689C19 #1047

BASE M/E: 91
RIC: 39487.

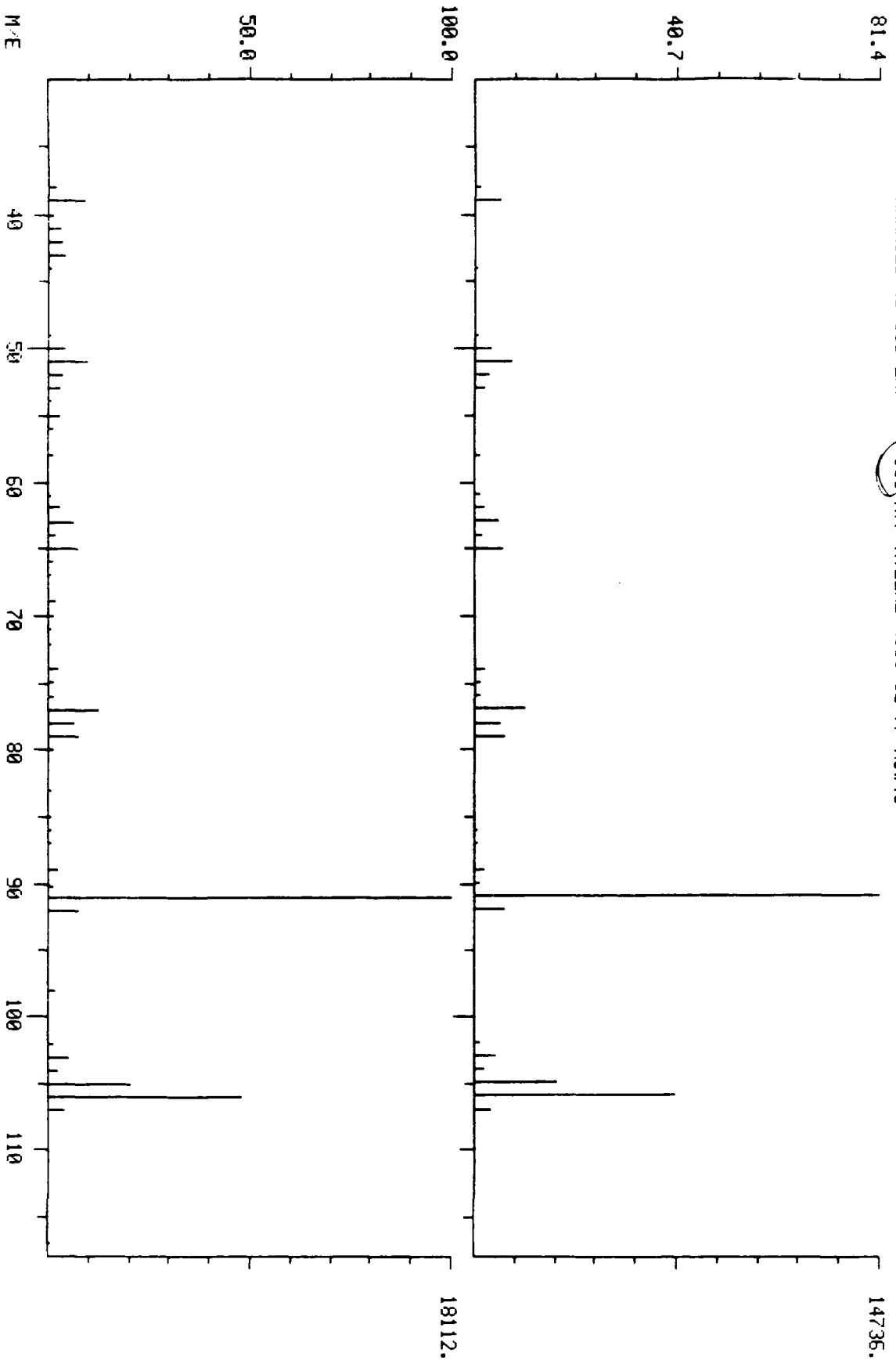
C8.H10
Y MT 1000
B PK 91
RANK 48
IN 48
PUR 931



COMPUCHEM LABS

DUAL MASS SPECTRUM
12/22/89 7:48:00 + 13:05
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B QM#19
ENHANCED (5 158 2N) 330 M,P-XYLENE <133-02-7> R0#48

DATA: GH009689C19 #1047 BASE M/E: 91/ 91
RIC: 39999.7 54207.

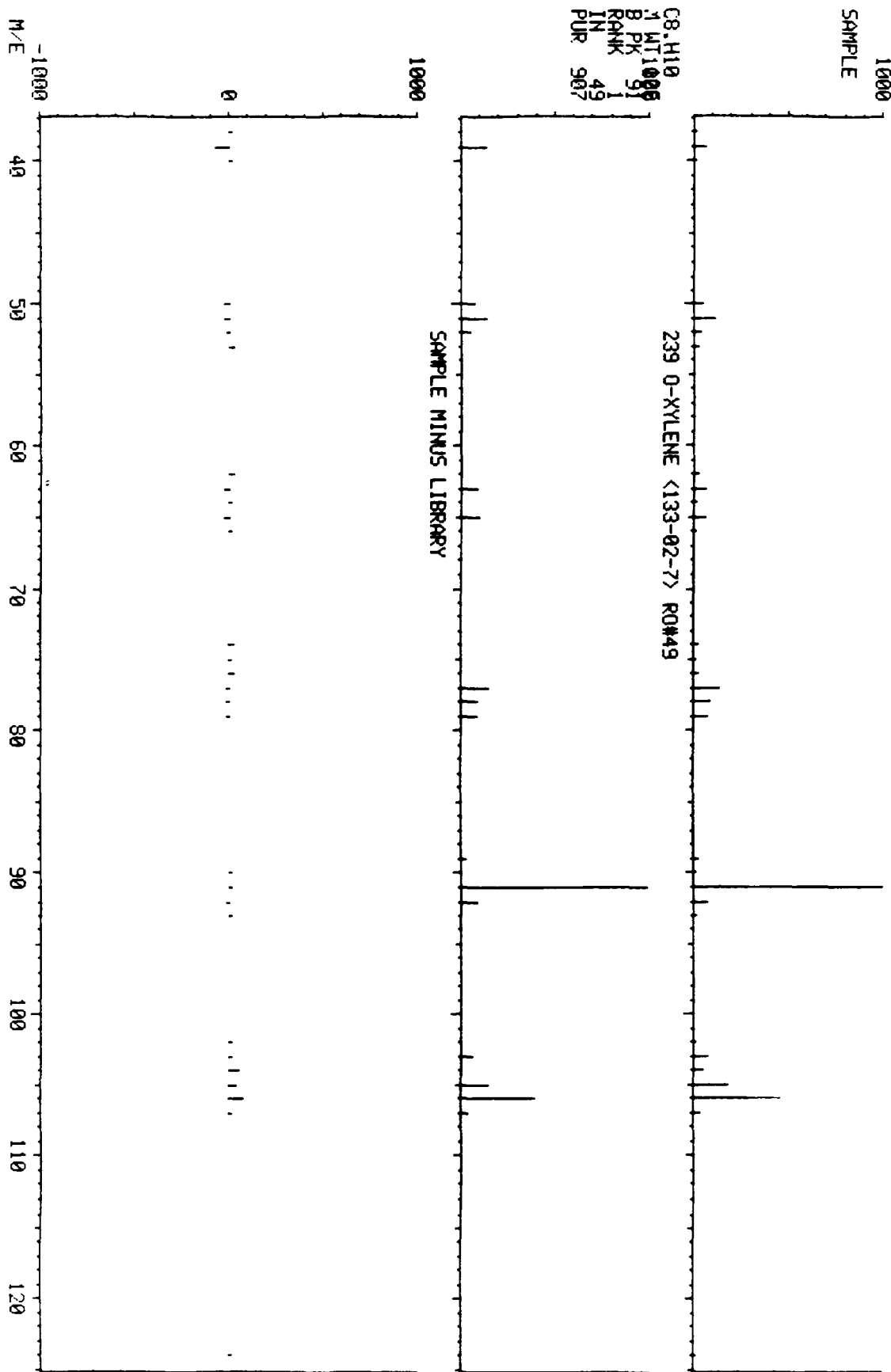


COMPUchem LABS
LIBRARY SEARCH
12/22/89 7:40:00 + 13:47
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B ON#19
ENHANCED (5 15B 2N 0T)

DATA: GH009689C19 #1103

BASE M/E: 91
RIC: 29471.

C8.H10
1 MT 1000
8 PK 91
RANK 49
IN 1
PUR 907

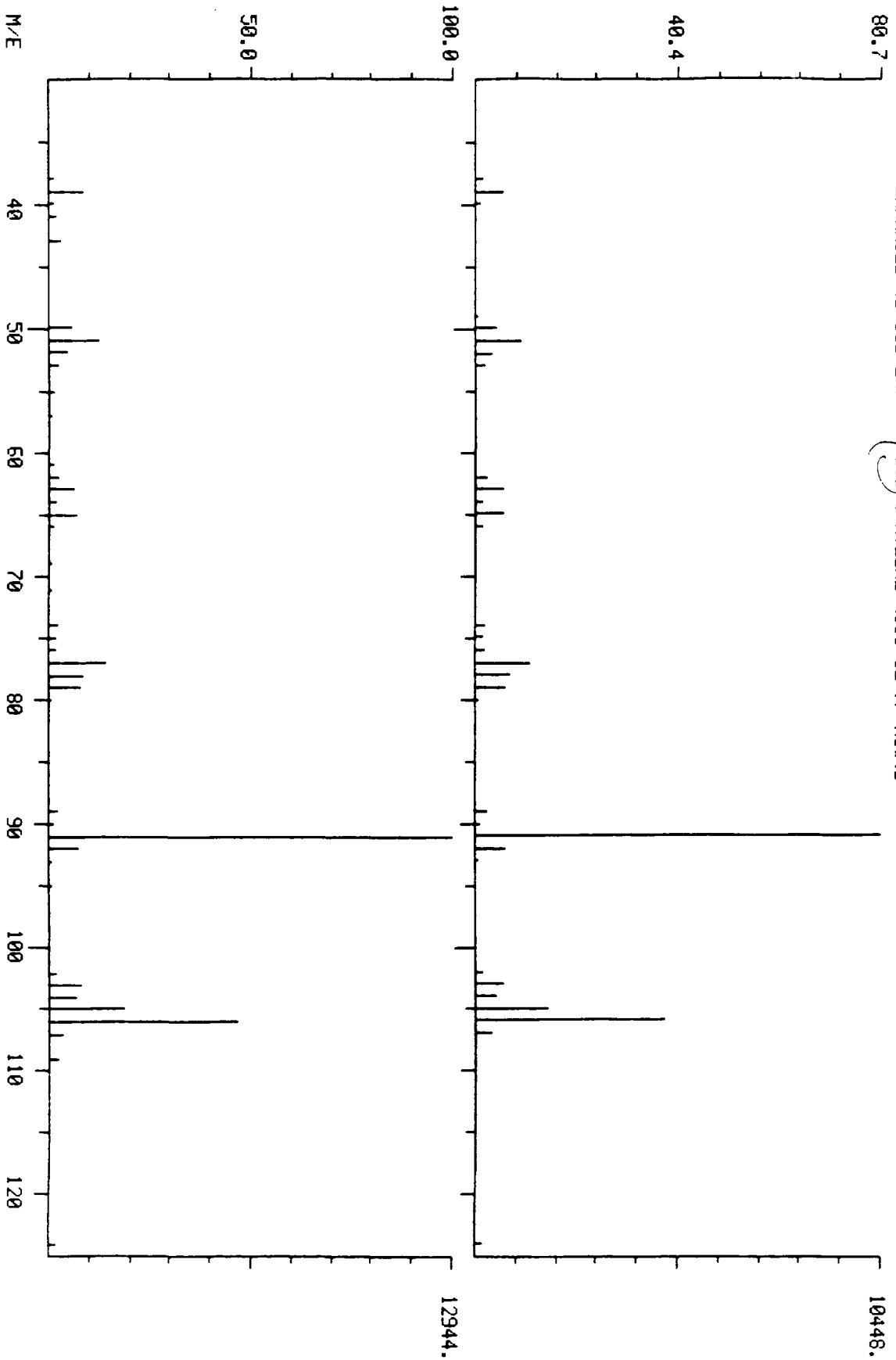


COMPUCHEM LABS

DATA: GH009689C19 #1103 BASE M/E: 91/ 91

RIC: 29695.7 32655.

DUAL MASS SPECTRUM
12/22/89 7:40:00 + 13:47
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (S 158 2N) 239 O-XYLENE <133-02-7> RO#49

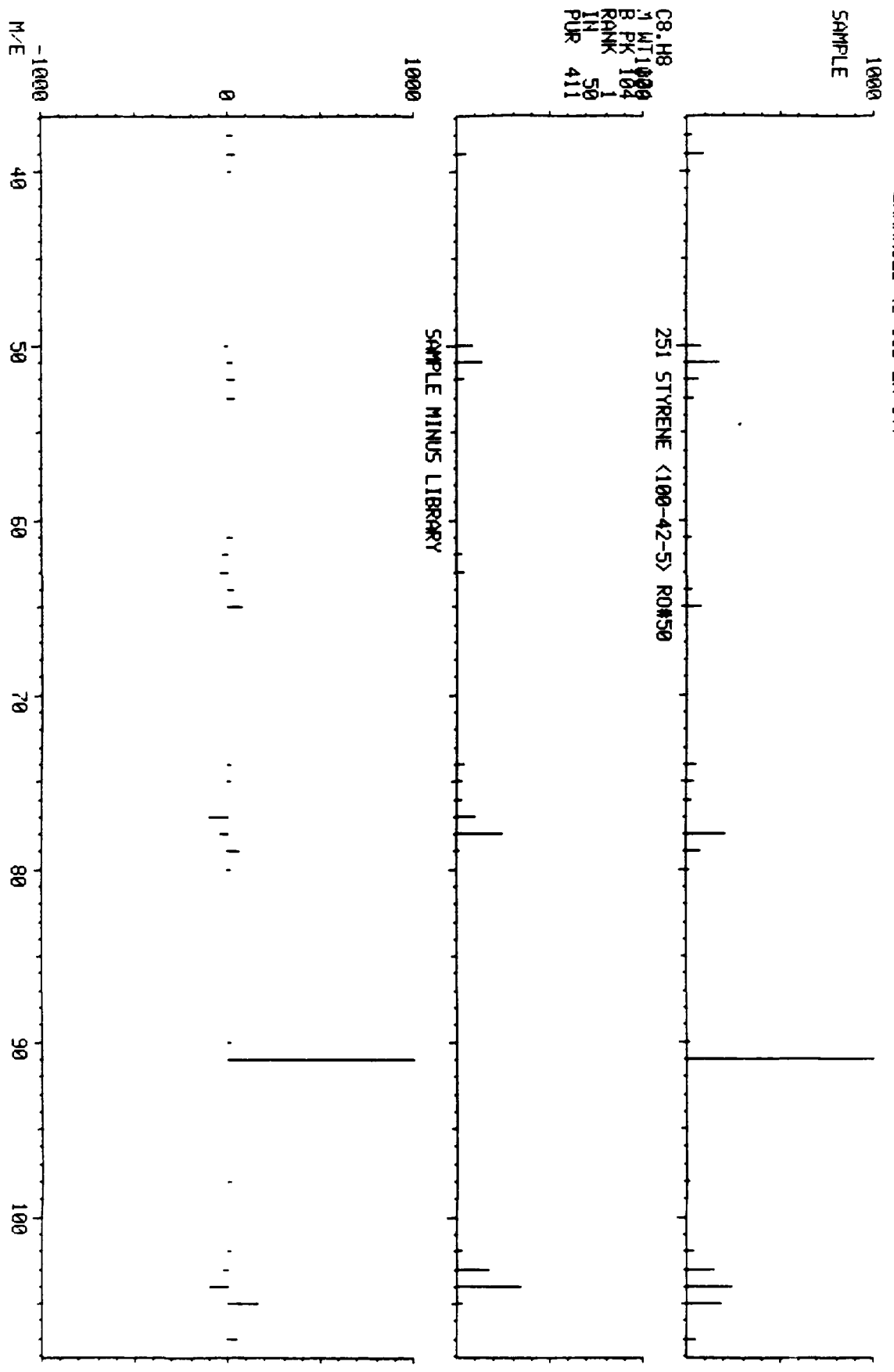


COMPUchem LABS
LIBRARY SEARCH
12/22/89 7:40:00 + 13:49
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#E202B ON#19
ENHANCED (5 158 2N 0T)

DATA: GH009689C19 #1106

BASE M/E: 91
RIC: 10063.

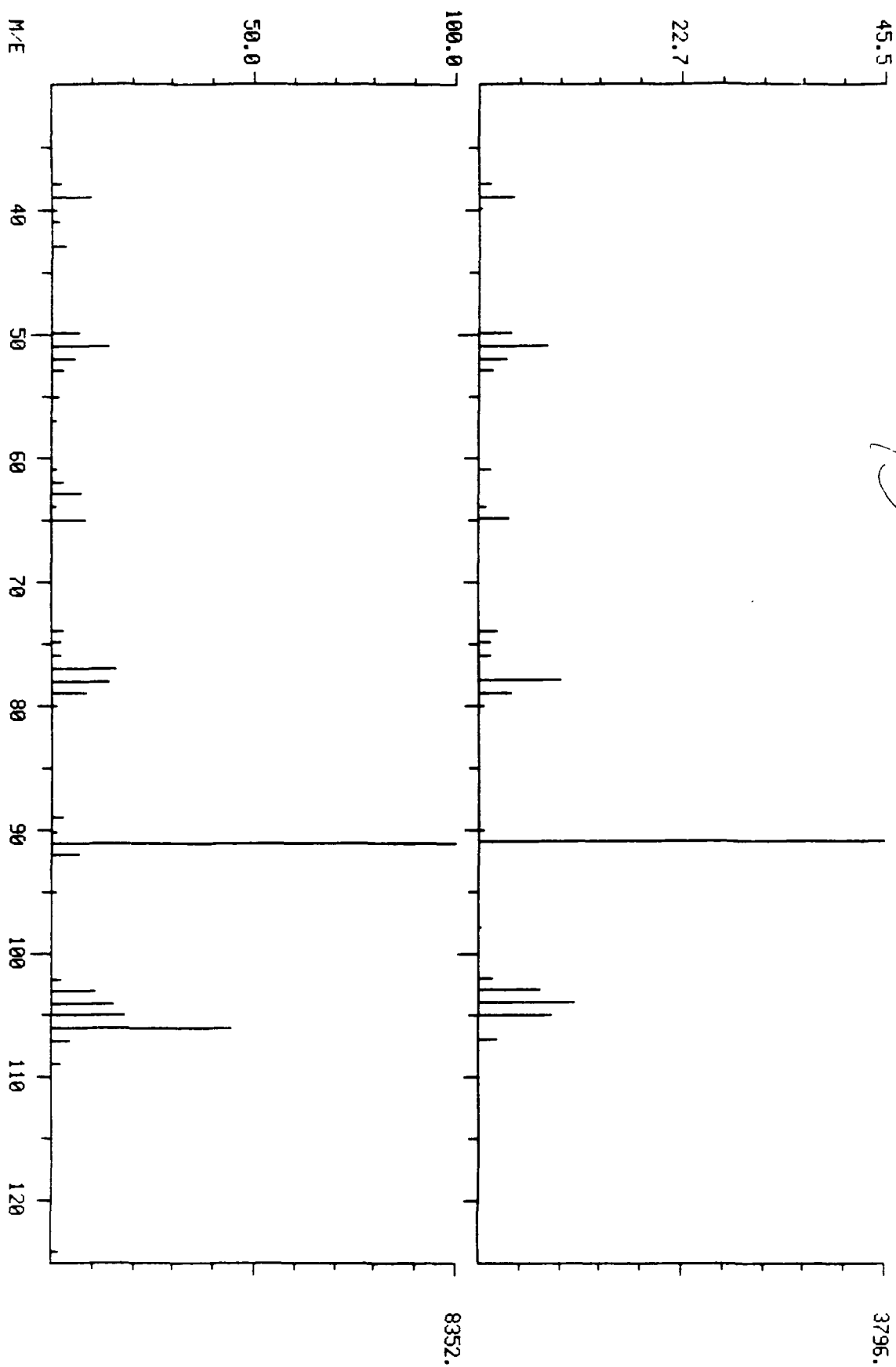
CG.H8
1 MT 1000
B PK 104
RANK 1
IN 50
PUR 411



DUAL MASS SPECTRUM
12/22/89 7:40:00 + 13:49
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B ON#19
ENHANCED (5 15B 2N) 251 STYRENE (100-42-5) RM#50

COMPUCHEN LABS

DATA: GH009689C19 #1106 BASE M/E: 91 / 91
RIC: 10063. / 27519.

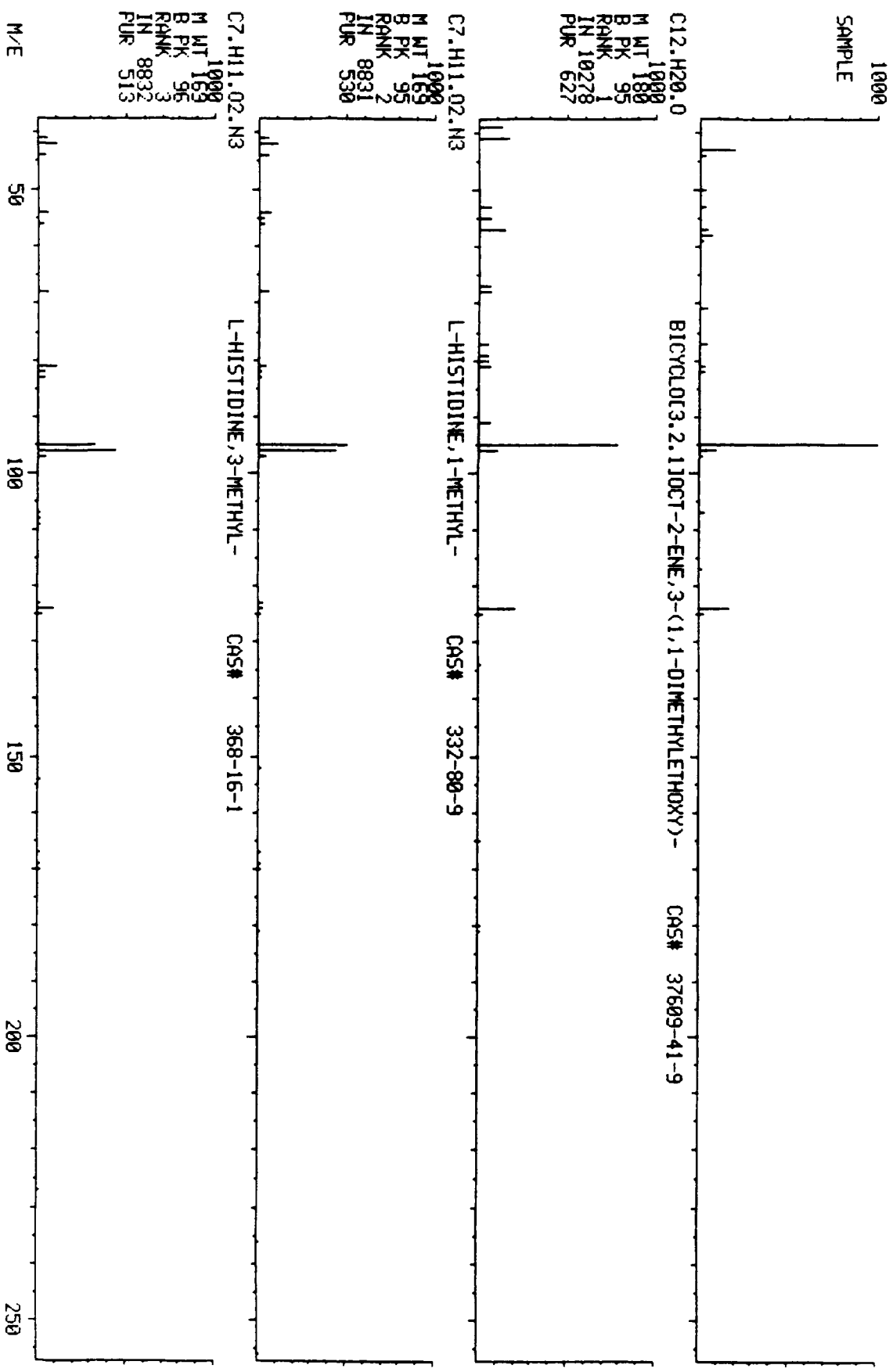


LIBRARY SEARCH
12/22/89 7:40:00 + 13:23
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (S 158 ZN 0T)

COMPUCHEM LABS

DATA: GH009689C19 #1071

BASE M/E: 95
RIC: 4887.

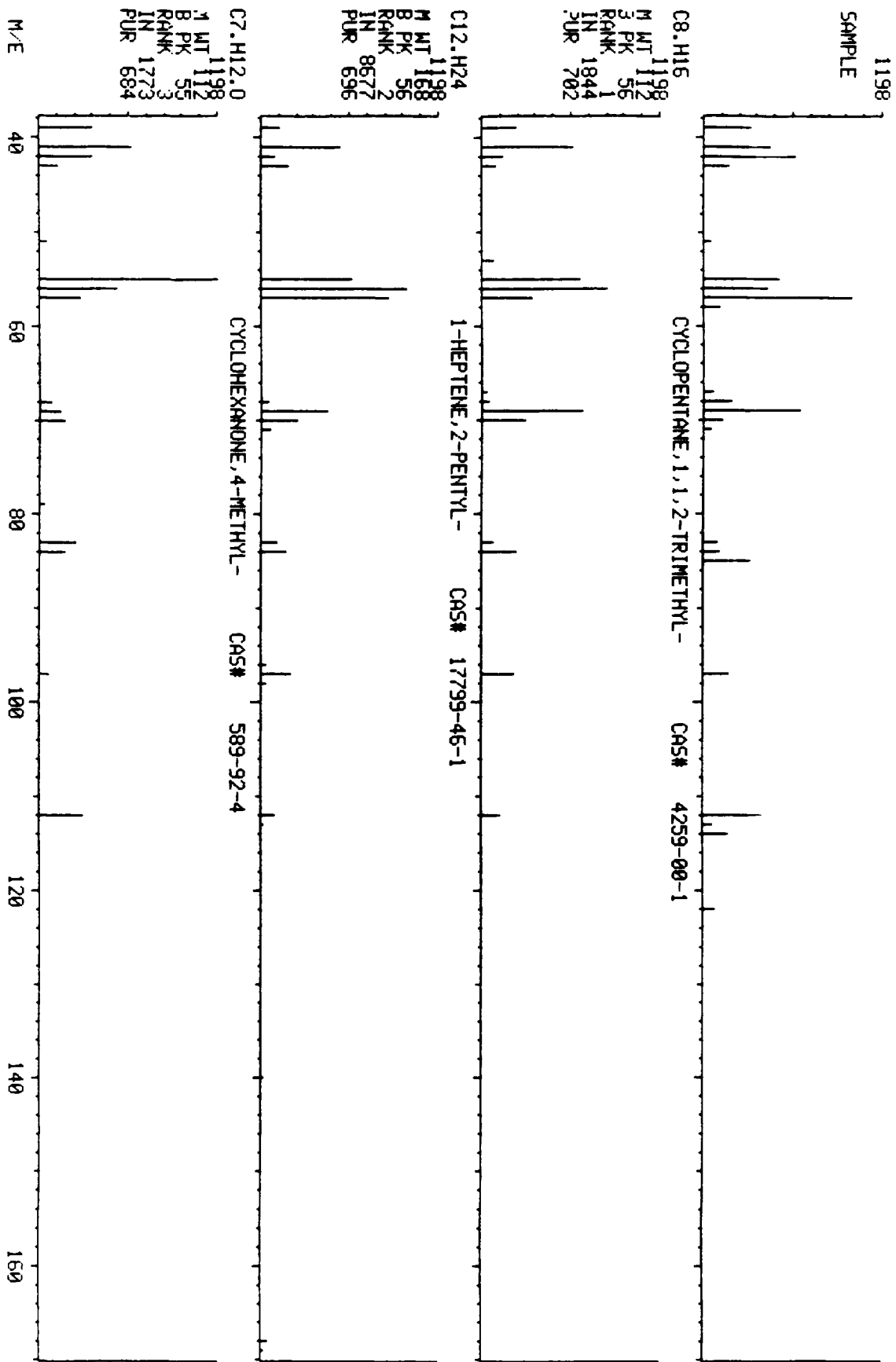


LIBRARY SEARCH
12/22/89 7:40:00 + 14:03
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B202B ON#19
ENHANCED (S 158 2N 0T)

COMPUCHEM LABS

DATA: GH009689C19 #1124

BASE M/E: 57
RIC: 7303.

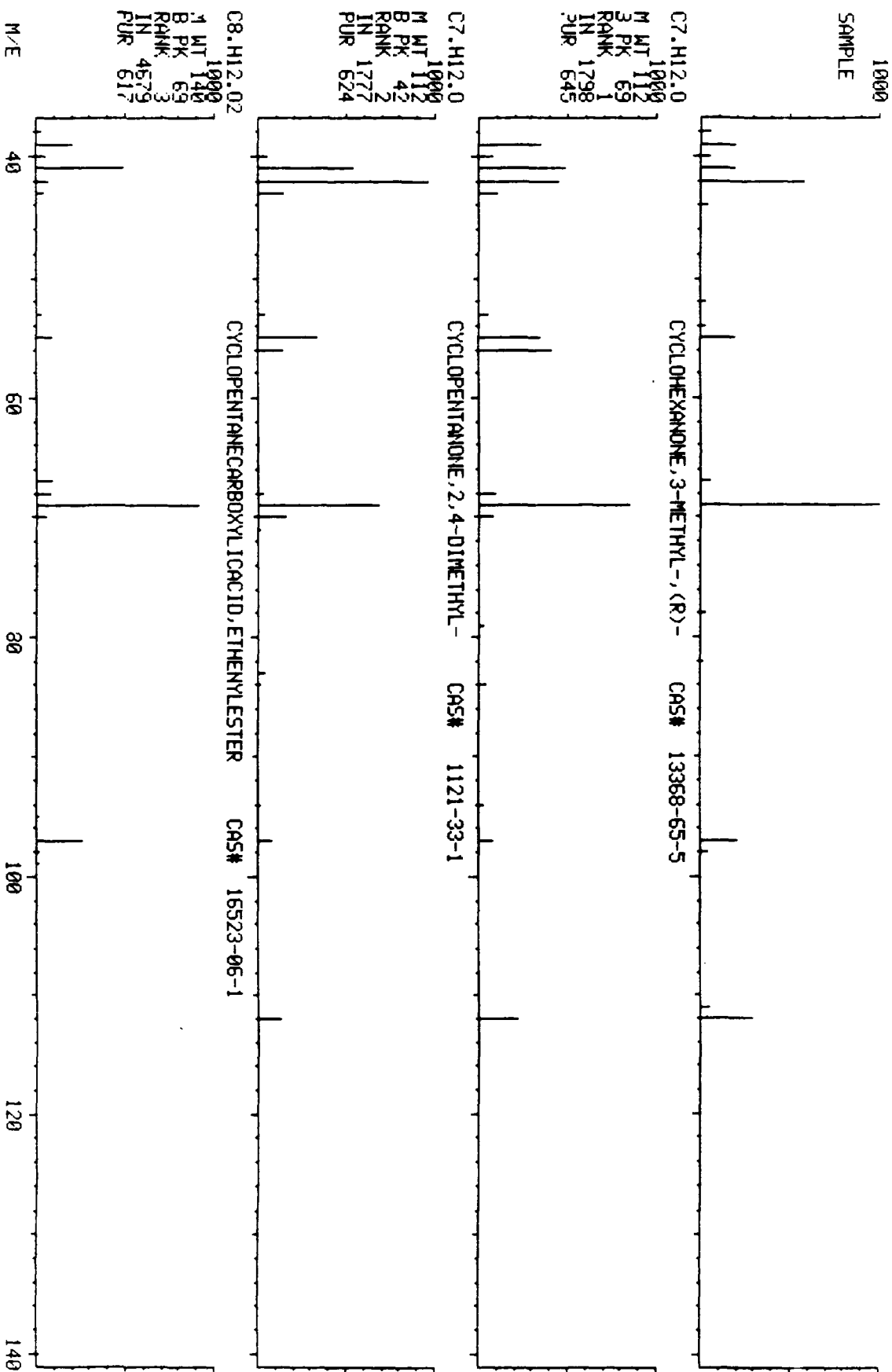


LIBRARY SEARCH
 12/22/89 7:40:00 + 14:24
 SAMPLE: 5G CC#309689 CASE#18756.7 EPARB2028 ON#19
 ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009689C19 #1152

BASE M/E: 59
 RIC: 6431.

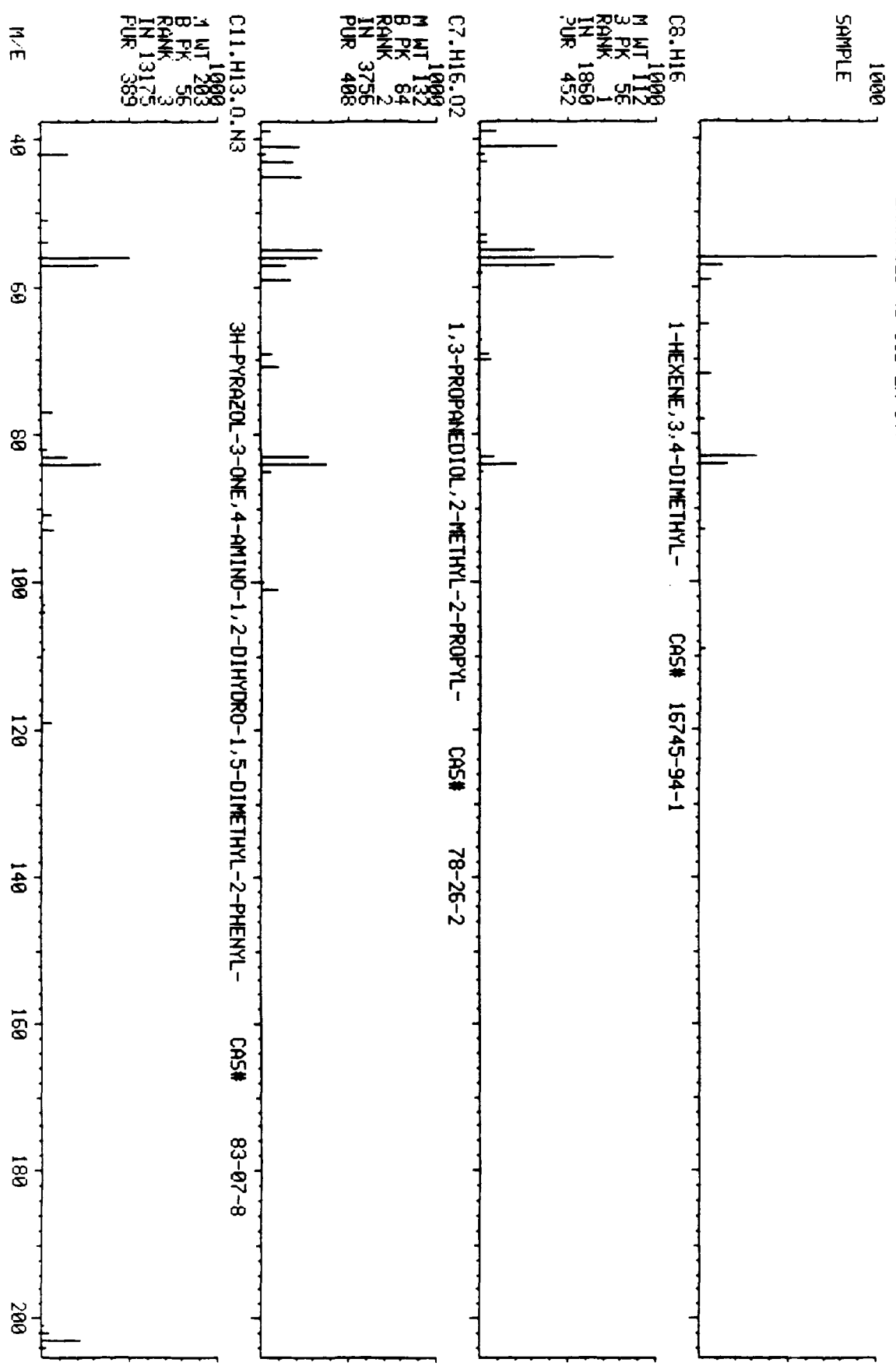


LIBRARY SEARCH
 12/22/89 7:40:00 + 14:29
 SAMPLE: SG CC#309689 CASE#18756.7 EPA#B202B ON#19
 ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009689C19 #1159

BASE M/E: 56
 RIC: 2447.



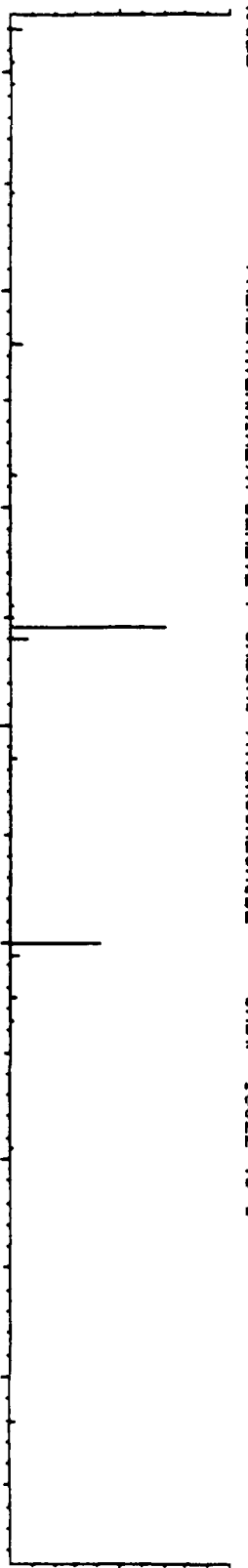
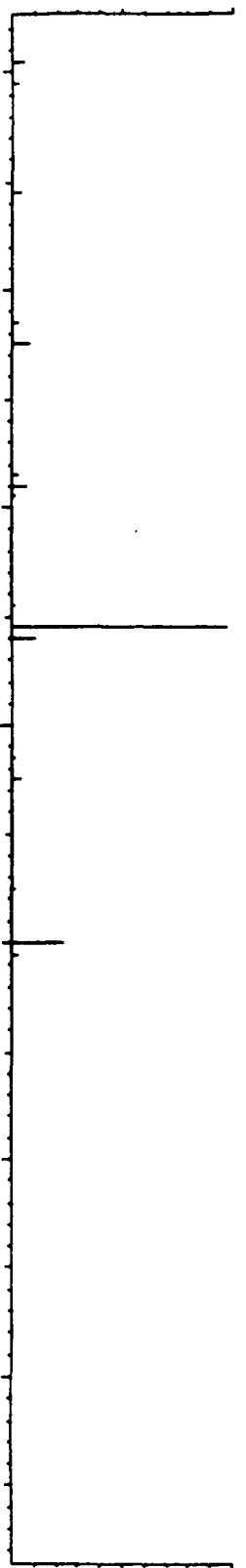
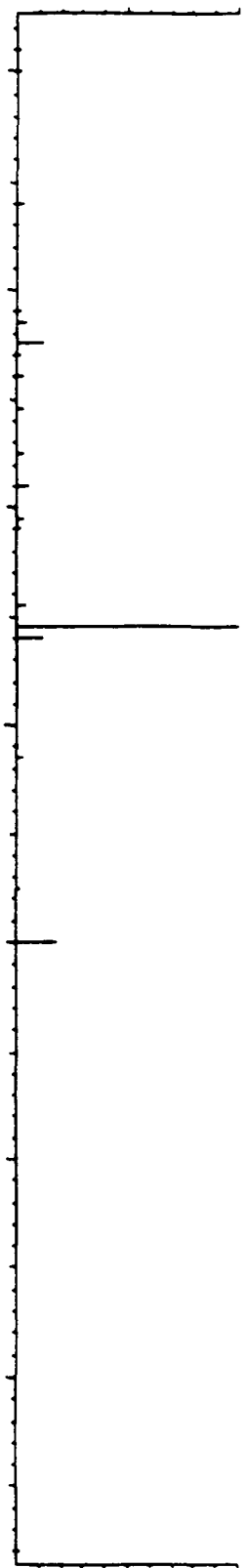
LIBRARY SEARCH
 12/22/89 7:46:00 + 15:21
 SAMPLE: 5G CCM309689 CASE#18756.7 EPA#B2028 QN#19
 ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

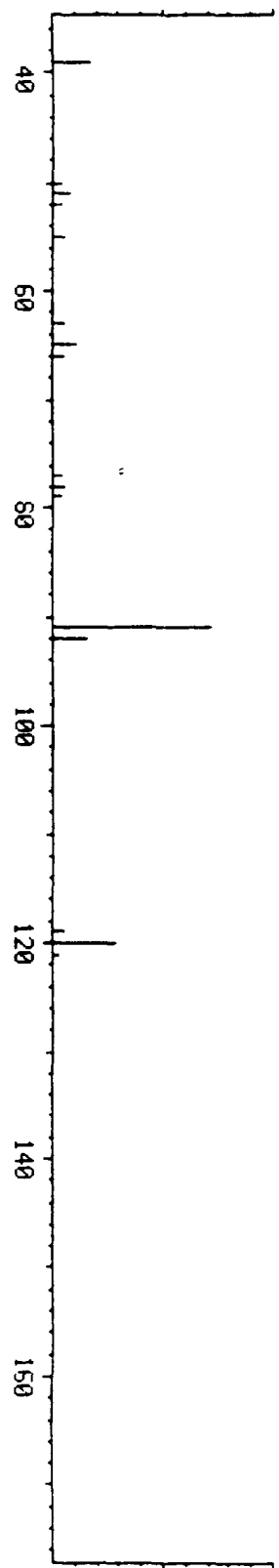
DATA: GH009689C19 #1228

BASE M/E: 91
 RIC: 4903.

1000
 SAMPLE



M/E

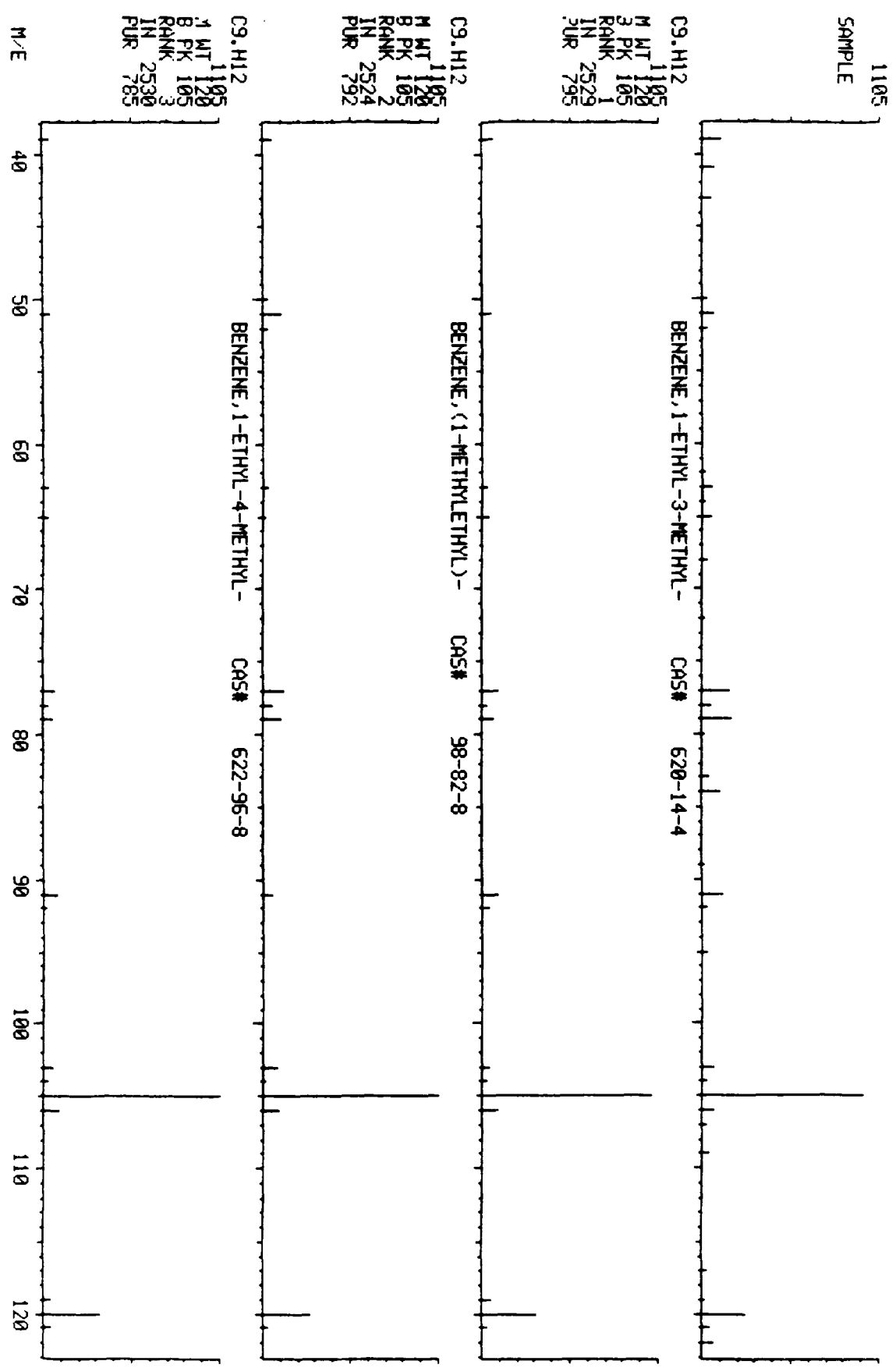


LIBRARY SEARCH
 12/22/89 7:40:00 + 15:34
 SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
 ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

DATA: GH009689C19 #1245

BASE M/E: 105
 RIC: 30879.



LIBRARY SEARCH
12/22/89 7:40:00 + 16:02
SAMPLE: 5G CC#309689 CASE#18756.7 EPA#B2028 ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: GH009689C19 #1283

BASE M/E: 105
RIC: 17055.

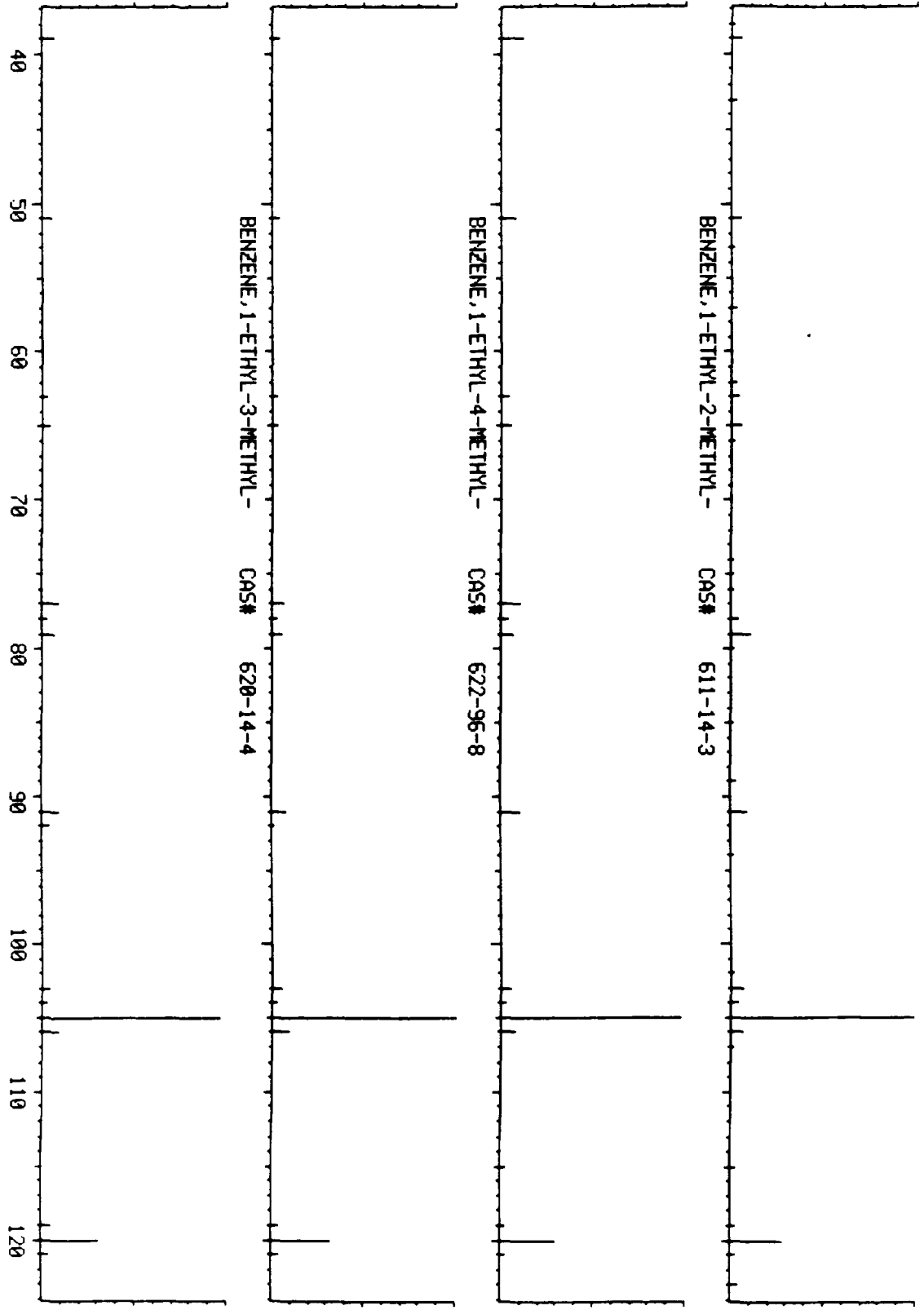
1014
SAMPLE

C9.H12
M WT 1014
PK 120
RANK 105
IN 2528
PUR 859

C9.H12
M WT 1014
PK 120
RANK 105
IN 2530
PUR 843

C9.H12
M WT 1014
PK 120
RANK 105
IN 2529
PUR 832

M/E



LAB INSTRUCTIONS:

RECEIPT DATE 12/20/89

CASE#: 18756 7

DUE DATE:

VOA
GC/MS WORKSHEET

COMPUCHEM#: 309689

R1 [] R2 [] D1 [] ()
R3 [] R4 [] D2 [] ()

L L SOLID, EPA SOW 2/88

Sample Prep Code---155
Instrument Code----413
Compound List-----494
Surrogate Std-----394
Internal Std-----036

=====

SAMPLE ID#: B202B Dry Wt. Factor 1.16 % Moisture 14

=====

GC/MS ANALYSIS

Amount Purged: [/] 10mls/Xg soil or [] Dilution _____ ul/10000ul Xg soil
Internal Standard Volume Added 5 ul
Surrogate Standard Volume Added 5 ul
BFB Filename B6891221.D19 Disk (29524)
Blank Filename GH009705.C19 Disk ()
Standard Filename G5891222.C19 Disk ()
Sample Filename GH009689.C19 Disk ()

DEC 28 1989
WORK UP 1452

ANALYST(S): Injection 1875-26

GC/MS REVIEW

CONDITION
CODE

OK

Entry Codes OK, EA, ES, SM, JS, SL, SH, JA, DA

Non-Entry Codes IM, IL, IH, SW, CT, CS, PC, NR
IF, LA, DI, CO, RN, DW, SI, SF
UP, BB, OT, VC, FO, NS

Extraneous Peak Search Results:

of Peaks Found: 07

[Handwritten signature]

Disposition: [/] Complete
[] Reprep neat required
[] Reprep using _____ g
[] Dilute ()

Quality Assurance Notice(s):

Notices Required 0

COMMENTS:

GC/MS Review OK Date 12/27/89 Auditor SKR Date 12/20/89

REPORT INTEGRATION

Final Reportable Package(s): G40-C19

Total # of Injections: 1

QA COMMENTS:

FINAL REVIEW:

Initials _____ Date ____/____/____

Initials _____ Date ____/____/____

AC1007 (05/89)

VOLATILE PREPARATION WORKSHEET

QUEUE # 19 PREP CODE -155 ASSIGNED TO Harshad Jeshi DATE 12/20/89

SAMPLE NUMBER	CASE NUMBER	SDG	QC SAMPLE		SAMPLE WEIGHT (G) VOLUME (ML)	SAMPLE ID	COMMENTS
			TYPE	ORIGINAL			
309763	18520	0164			5.0g	TB-12	
309768	↓	↓			5.0g	WB5-1	
309769	↓	↓			5.0g	WB5-2	
309770	↓	↓			5.0g	WB5-4	
309679	18756	0007			5.0g	B201A	
309680	↓	↓	SS	309679	5.0g	B201A MS	
309681	↓	↓	SS	309679	5.0g	B201A MSD	
309682	↓	↓	BS		0.0g		
309686	↓	↓			5.0g	B201B	
309687	↓	↓			5.0g	B202AR	
309688	↓	↓			5.0g	B202A	
309689	↓	↓			5.0g	B202B	
309690	↓	↓			5.0g	B202C	
309784			B1		5.0 ml	B3	
309785			B2		0.0 ml	B4	
309786			B3		0.0 ml	B5	
309787			B4		0.0 ml	B6	
309788			B5		0.0 ml	B7	

SURROGATE # _____ LOT # _____ MANUAL OPERATOR 739 / 516
 AMOUNT 4
 RELINQUISHED BY [Signature] DATE 12/20/89 RECEIVED BY [Signature] DATE 12/20

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT. LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	465	46800	50.0		X-16
221	50	CHLOROMETHANE				BDL	12.15
231	62	VINYL CHLORIDE				BDL	1
220	94	BROMOMETHANE				BDL	1
209	64	CHLOROETHANE				BDL	1
216	96	1,1-DICHLOROETHENE				BDL	1
254	76	CARBON DISULFIDE				BDL	
252	43	ACETONE (2-PROPANONE)			45.6	46.5313	1
248	114 I	1,4-DIFLUOROBENZENE (IS)	609	192000	50.0		
222	84	METHYLENE CHLORIDE			26.0	26.3012	
226	96	TRANS-1,2-DICHLOROETHENE				BDL	
214	63	1,1-DICHLOROETHANE				BDL	
257	43	VINYL ACETATE				BDL	1
237	96	CIS-1,2-DICHLOROETHENE				BDL	
253	72	2-BUTANONE			34.1	34.40	1
211	83	CHLOROFORM				BDL	
227	97	1,1,1-TRICHLOROETHANE				BDL	
206	117	CARBON TETRACHLORIDE				BDL	
203	78	BENZENE				BDL	
215	62	1,2-DICHLOROETHANE				BDL	
270	117 I	D5-CHLOROBENZENE (IS) RO#29	1000	204000	50.0		
229	130	TRICHLOROETHENE				BDL	
217	63	1,2-DICHLOROPROPANE				BDL	5
212	83	BROMODICHLOROMETHANE				BDL	5
218	75	CIS-1,3-DICHLOROPROPENE				BDL	5
256	43	4-METHYL-2-PENTANONE			2.4	BDL 2.4	10
225	92	TOLUENE			12.8	13.15	5
250	75	TRANS-1,3-DICHLOROPROPENE				BDL	5
228	97	1,1,2-TRICHLOROETHANE				BDL	5
224	164	TETRACHLOROETHENE				BDL	5
255	43	2-HEXANONE			20.2	20.23	10
208	129	DIBROMOCHLOROMETHANE				BDL	5
207	112	CHLOROBENZENE				BDL	5
219	106	ETHYLBENZENE			10.1	10.12	5
330	106	M, P-XYLENE			25.8	26.30	5
239	106	O-XYLENE			21.0	21.24	5
251	104	STYRENE			2.7	2.735	5
205	173	BROMOFORM				BDL	5
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	5
258	65 S	D4-1,2-DICHLOROETHANE RO#57			48.3	97.0%	
247	95 S	BROMOFLUOROBENZENE			47.8	96.0%	
233	98 S	D8-TOLUENE RO#59			45.8	92.0%	
289	106	XYLENES (TOTAL)			46.8	47.54	5

CORRECTED/REVIEWED BY OKA
(GC/MS DATA REVIEWER)

DATE 12.27.89

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT. LIMIT (UG/KG)
299	96	1,2-DICHLOROETHENE (TOTAL)				BDL	
CHECKSUMS:							
	3979.		2074	442800.		539.4	533.

CORRECTED/REVIEWED BY *OK*
(GC/MS DATA REVIEWER)

DATE 12-27-89

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P	F
40	258	D4-1,2-DICHLOROETHANE RO#57	48.3	50.0	97.	70-121	X	
41	247	BROMOFLUOROBENZENE	47.8	50.0	96.	74-121	X	
42	233	D8-TOLUENE RO#59	45.8	50.0	92.	81-117	X	

* ADVISORY SURROGATE ONLY

++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

=====

CORRECTION FACTOR CALCULATION:

$$\frac{5.0 \text{ G}}{\text{WET WEIGHT OF SAMPLE (G)}} \times \frac{\text{GC/MS DILUTION FACTOR}}{\text{GC/MS DILUTION FACTOR}} \times \frac{\text{DRY WEIGHT FACTOR}}{\text{DRY WEIGHT FACTOR}} =$$

$$\frac{5.0 \text{ G}}{5.00 \text{ (G)}} \times \frac{1.0}{1.00} \times \frac{1.16}{1.16} = 1.000$$

=====

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:

THE SURROGATES ARE ADDED TO THE SAMPLE PRIOR TO SPARGING.
SURROGATE SPIKE CONVERSION FACTOR = 1.

=====

VERSION 8

CORRECTED/REVIEWED BY P. H. L.
(GC/MS DATA REVIEWER)

DATE 12 27 89

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202C

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309690
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009690A19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 13 Date Analyzed: 12/22/89
 Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	11	U
74-83-9	-----Bromomethane	11	U
75-01-4	-----Vinyl Chloride	11	U
75-00-3	-----Chloroethane	11	U
75-09-2	-----Methylene Chloride	23	B
67-64-1	-----Acetone	26	B
75-15-0	-----Carbon Disulfide	6	U
75-35-4	-----1,1-Dichloroethene	6	U
75-34-3	-----1,1-Dichloroethane	6	U
540-59-0	-----1,2-Dichloroethene (total)	6	U
67-66-3	-----Chloroform	6	U
107-06-2	-----1,2-Dichloroethane	6	U
78-93-3	-----2-Butanone	3	J
71-55-6	-----1,1,1-Trichloroethane	6	U
56-23-5	-----Carbon Tetrachloride	6	U
108-05-4	-----Vinyl Acetate	11	U
75-27-4	-----Bromodichloromethane	6	U
78-87-5	-----1,2-Dichloropropane	6	U
10061-01-5	-----cis-1,3-Dichloropropene	6	U
79-01-6	-----Trichloroethene	6	U
124-48-1	-----Dibromochloromethane	6	U
79-00-5	-----1,1,2-Trichloroethane	6	U
71-43-2	-----Benzene	6	U
10061-02-6	-----Trans-1,3-Dichloropropene	6	U
75-25-2	-----Bromoform	6	U
108-10-1	-----4-Methyl-2-Pentanone	11	U
591-78-6	-----2-Hexanone	11	U
127-18-4	-----Tetrachloroethene	6	U
79-34-5	-----1,1,2,2-Tetrachloroethane	6	U
108-88-3	-----Toluene	2	J
108-90-7	-----Chlorobenzene	6	U
100-41-4	-----Ethylbenzene	1	J
100-42-5	-----Styrene	6	U
1330-20-7	-----Total Xylenes	7	

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202C

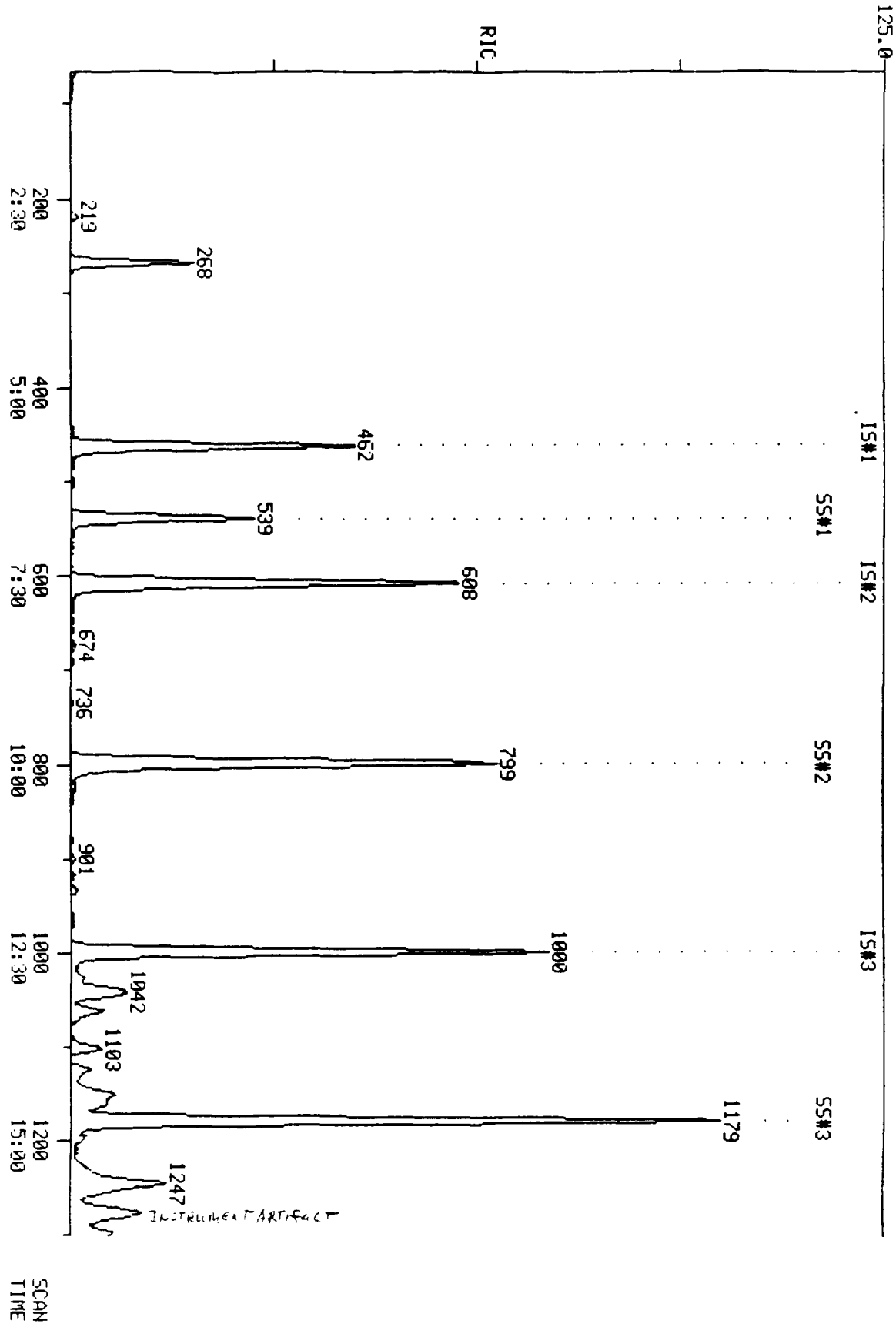
Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309690
 Sample wt/vol: 5.0 (g/mL) G Lab File ID: GH009690A19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 13 Date Analyzed: 12/22/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 3 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	13.02	9.2	J
2. 4971-18-0	CYCLOPENTANONE, 2-ETHYL-	15.59	14	J
3.	INSTRUMENT ARTIFACT	15.97	9.2	J

RIC
 12/22/89 8:15:00
 SAMPLE: 5 GRAMS CC#309690 EPA#:B202C CASE#18756.7 ON#19
 COND.:

COMPUCHEM LABS
 COMPUCHEM DATA: CH009690A19 SCANS 67 TO 1300



QUANTITATION REPORT FILE: GH009690A19
 DATA: GH009690A19.TI ✓
 12/22/89 8:15:00 ✓
 SAMPLE: 5 GRAMS CC#309690 EPA#: B202C CASE#18756.7 DN#19
 CONDS.:
 SUBMITTED BY: 19 ANALYST: 1492 ✓

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
 RESP. FAC. FROM LIBRARY ENTRY ✓

- NO NAME
- 1 *234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1
- 2 221 CHLOROMETHANE <74-87-3> RO#2
- 3 231 VINYL CHLORIDE <75-01-4> RO#3
- 4 220 BROMOMETHANE <78-83-9> RO#4
- 5 209 CHLOROETHANE <75-00-3> RO#5
- 6 216 1,1-DICHLOROETHENE <75-35-4> RO#8
- 7 254 CARBON DISULFIDE <75-15-0> RO#9
- 8 252 ACETONE (2-PROPANONE) <67-64-1> RO#13
- 9 *248 1,4-DIFLUOROBENZENE (IS) <540-36-3> RO#14
- 10 222 METHYLENE CHLORIDE <75-09-2> RO#16
- 11 226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17
- 12 214 1,1-DICHLOROETHANE <75-34-3> RO#19
- 13 257 VINYL ACETATE <108-05-4> RO#20
- 14 237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21
- 15 253 2-BUTANONE <78-93-3> RO#22
- 16 211 CHLOROFORM <67-66-2> RO#23
- 17 227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24
- 18 206 CARBON TETRACHLORIDE <56-23-5> RO#25
- 19 203 BENZENE <71-43-2> RO#26
- 20 215 1,2-DICHLOROETHANE <107-06-2> RO#27
- 21 *270 D5-CHLOROBENZENE (IS) RO#29
- 22 229 TRICHLOROETHENE <79-01-6> RO#30
- 23 217 1,2-DICHLOROPROPANE <78-87-5> RO#31
- 24 212 BROMODICHLOROMETHANE <75-27-4> RO#33
- 25 218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35
- 26 256 4-METHYL-2-PENTANONE <108-01-1> RO#36
- 27 225 TOLUENE <108-88-3> RO#37
- 28 250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38
- 29 228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39
- 30 224 TETRACHLOROETHENE <127-18-4> RO#41
- 31 255 2-HEXANONE <591-78-6> RO#42
- 32 208 DIBROMOCHLOROMETHANE <124-48-1> RO#43
- 33 207 CHLOROBENZENE <108-90-7> RO#45
- 34 219 ETHYLBENZENE <100-41-4> RO#47
- 35 330 M,P-XYLENE <133-02-7> RO#48
- 36 239 O-XYLENE <133-02-7> RO#49
- 37 251 STYRENE <100-42-5> RO#50
- 38 205 BROMOFORM <75-25-2> RO#51
- 39 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54
- 40 *258 D4-1,2-DICHLOROETHANE RO#57
- 41 *247 BROMOFLUOROBENZENE <460-00-4> RO#58
- 42 *233 D8-TOLUENE RO#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	462	5:46	1	1.000	A BV	49081.	50.000 UG/KG	14.22
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	219	2:44	1	0.474	A BB	5922.	22.799 UG/KG	6.49%
9	114	608	7:36	9	1.000	A BB	202282.	50.000 UG/KG	14.22
10	84	268	3:21	1	0.580	A BB	31695.	19.649 UG/KG	5.59%
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	446	5:34	1	0.965	A BB	317.	3.013 UG/KG	0.86%
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	NOT FOUND							
20	62	NOT FOUND							
21	117	999	12:29	21	1.000	A BB	211350.	50.000 UG/KG	14.22
22	130	NOT FOUND							
23	63	NOT FOUND							
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	NOT FOUND							
27	92	809	10:07	21	0.810	A BB	5280.	1.692 UG/KG	0.48%
28	75	NOT FOUND							
29	97	NOT FOUND							
30	164	NOT FOUND							
31	43	920	11:30	21	0.921	A BB	2639.	3.198 UG/KG	0.91%
32	129	NOT FOUND							
33	112	NOT FOUND							
34	106	1048	13:06	21	1.049	A BB	2639518.	4.443 UG/KG	1.26%
35	106	1048	13:06	21	1.049	A BB	9518.	3.196 UG/KG	0.91%
36	106	1103	13:47	21	1.104	A BB	7427.	2.825 UG/KG	0.80%
37	104	NOT FOUND							
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	538	6:43	1	1.165	A BB	98864.	48.267 UG/KG	13.73
41	95	1179	14:44	21	1.180	A BB	164284.	45.433 UG/KG	12.93
42	98	799	9:59	21	0.800	A BB	208117.	46.995 UG/KG	13.37

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:54	0.98	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:06		10.000			50.00		0.952	
3	1:11		10.000			50.00		1.256	
4	1:26		10.000			50.00		1.579	
5	1:34		10.000			50.00		0.796	
6	2:37		5.000			50.00		1.346	
7	2:46		5.000			50.00		3.709	
8	2:48	0.98	10.000	0.05	22.80	50.00	0.121	0.265	0.46
9	7:42	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	3:27	0.97	5.000	0.12	19.65	50.00	0.646	1.643	0.39
11	3:53		5.000			50.00		1.486	
12	4:34		5.000			50.00		2.635	
13	4:52		10.000			50.00		0.458	
14	5:33		5.000			50.00		1.675	

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:42	0.98	10.000	0.10	3.01	50.00	0.006	0.107	0.06
16	6:10		5.000			50.00		3.433	
17	6:18		5.000			50.00		0.814	
18	6:34		5.000			50.00		0.755	
19	6:53		5.000			50.00		0.895	
20	6:58		5.000			50.00		2.523	
21	12:34	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	8:01		5.000			50.00		0.569	
23	8:22		5.000			50.00		0.352	
24	8:55		5.000			50.00		0.701	
25	9:41		5.000			50.00		0.690	
26	10:04		15.000			50.00		0.257	
27	10:10	0.99	5.000	0.16	1.69	50.00	0.025	0.738	0.03
28	10:42		5.000			50.00		0.342	
29	10:58		5.000			50.00		0.349	
30	11:06		5.000			50.00		0.723	
31	11:33	1.00	15.000	0.06	3.20	50.00	0.012	0.195	0.06
32	11:35		5.000			50.00		0.690	
33	12:37		5.000			50.00		1.092	
34	12:55	1.01	5.000	0.21	4.44	50.00	0.045	0.507	0.09
35	13:10	1.00	5.000	0.21	3.20	50.00	0.045	0.705	0.06
36	13:52	0.99	5.000	0.22	2.83	50.00	0.035	0.622	0.06
37	13:54		5.000			50.00		1.099	
38	14:07		5.000			50.00		0.664	
39	15:16		5.000			50.00		0.557	
40	6:52	0.98	5.000	0.23	48.27	50.00	2.014	2.087	0.97
41	14:49	0.99	5.000	0.24	45.43	50.00	0.777	0.855	0.91
42	10:04	0.99	5.000	0.16	46.99	50.00	0.985	1.048	0.94

LIBRARY SEARCH
12/22/89 8:15:00 + 2:44
SAMPLE: 5 GRAMS CC#309690 EPA#:B202C CASE#18756.7 ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: CH009690R19 # 219

BASE M/E: 43
RIC: 938.

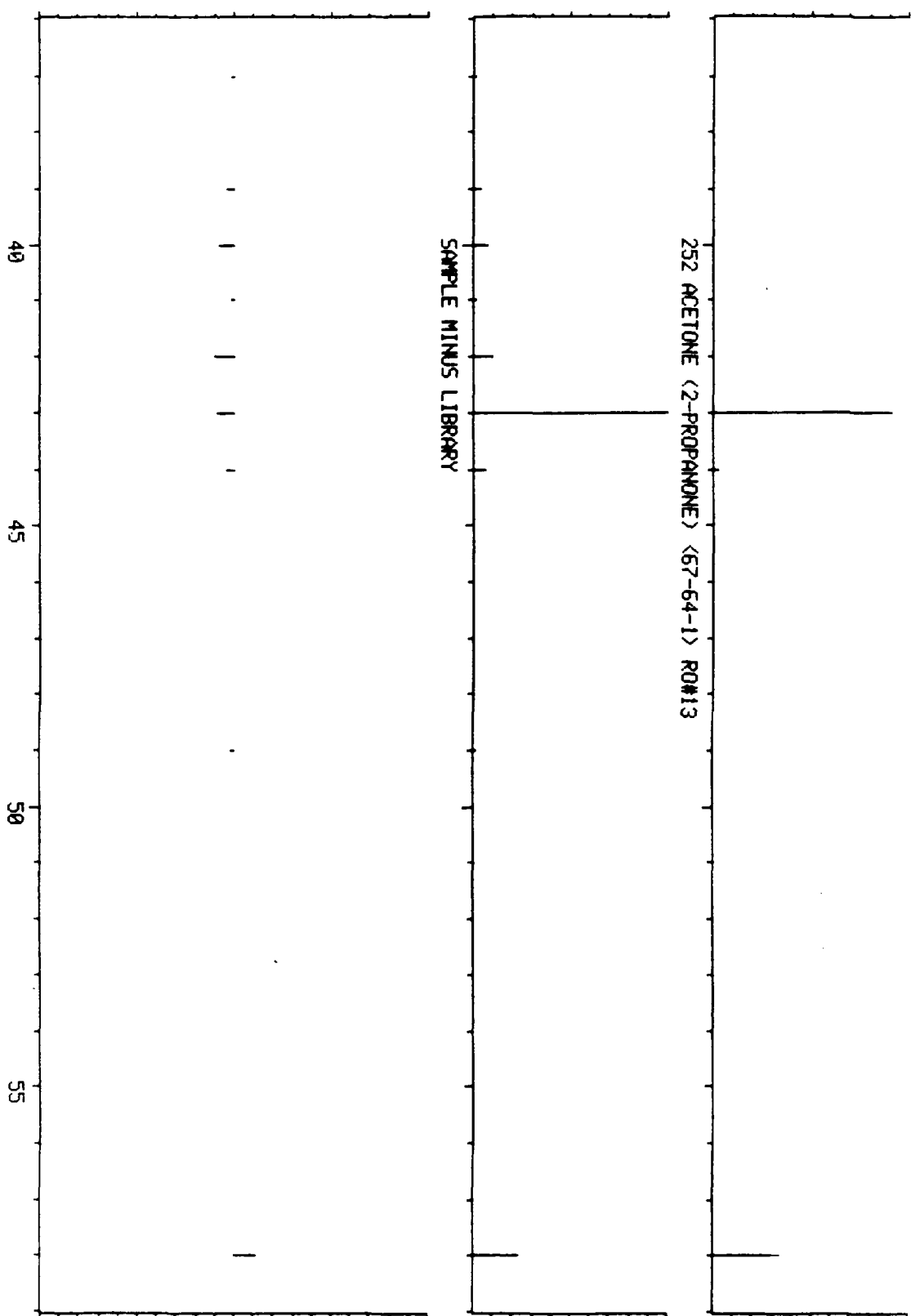
1092
SAMPLE

C3.H6.O
M WT 1092
B PK 43
RANK 1
IN 13
PUR 846

252 ACETONE (2-PROPANONE) (67-64-1) R0#13

SAMPLE MINUS LIBRARY

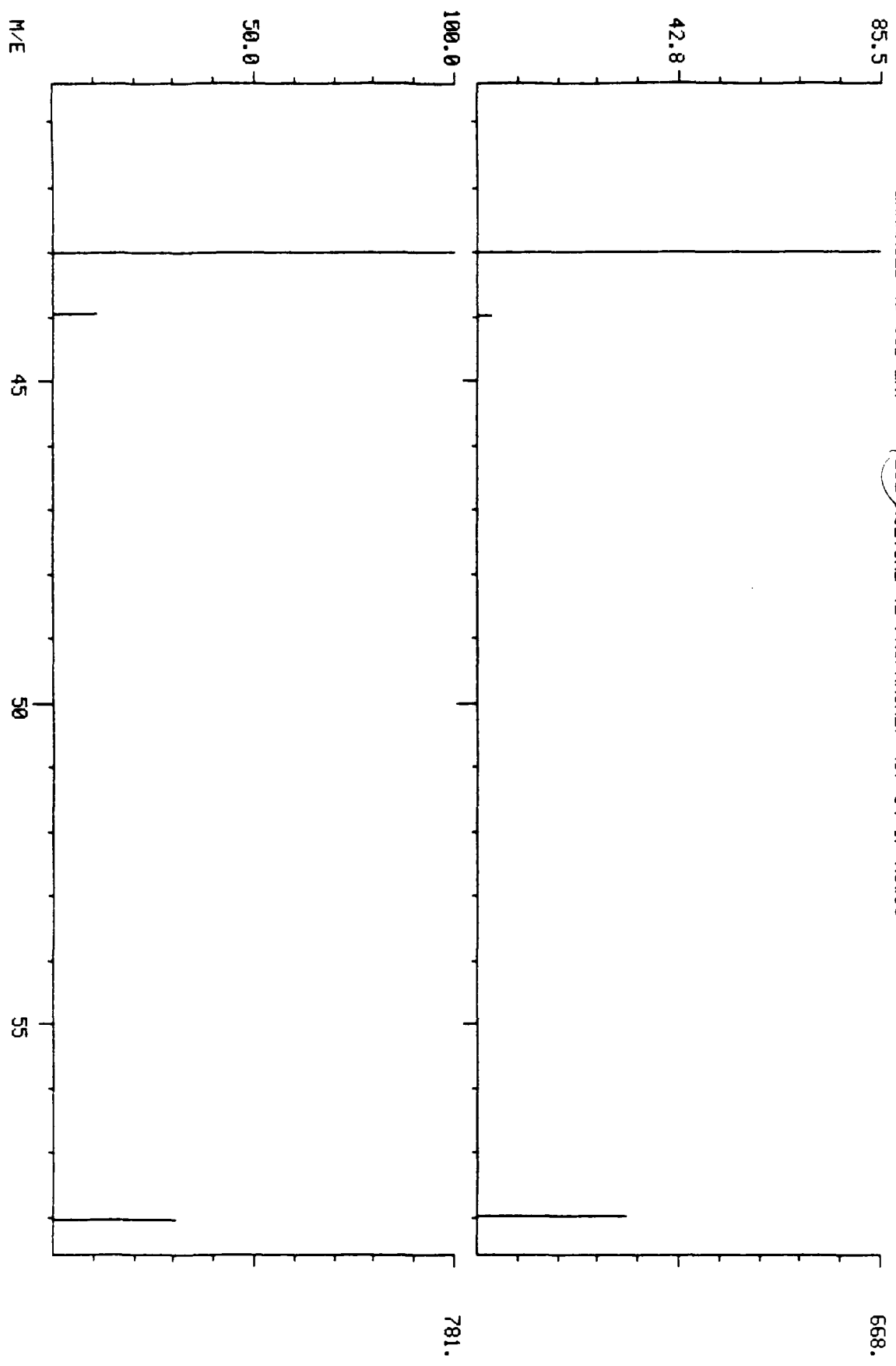
-1092
M/E



DUAL MASS SPECTRUM
12/22/89 8:15:00 + 2:44
SAMPLE: 5 GRAMS OC#309690 EPA#: B202C CASE#18756.7 ON#19
ENHANCED (5 15B 2N) (252 ACETONE (2-PROPANONE) <67-64-1> R0#13

COMPUCHEM LABS

DATA: GH009690A19 #219 BASE M/E: 43/ 43
RIC: 938./ 1105.



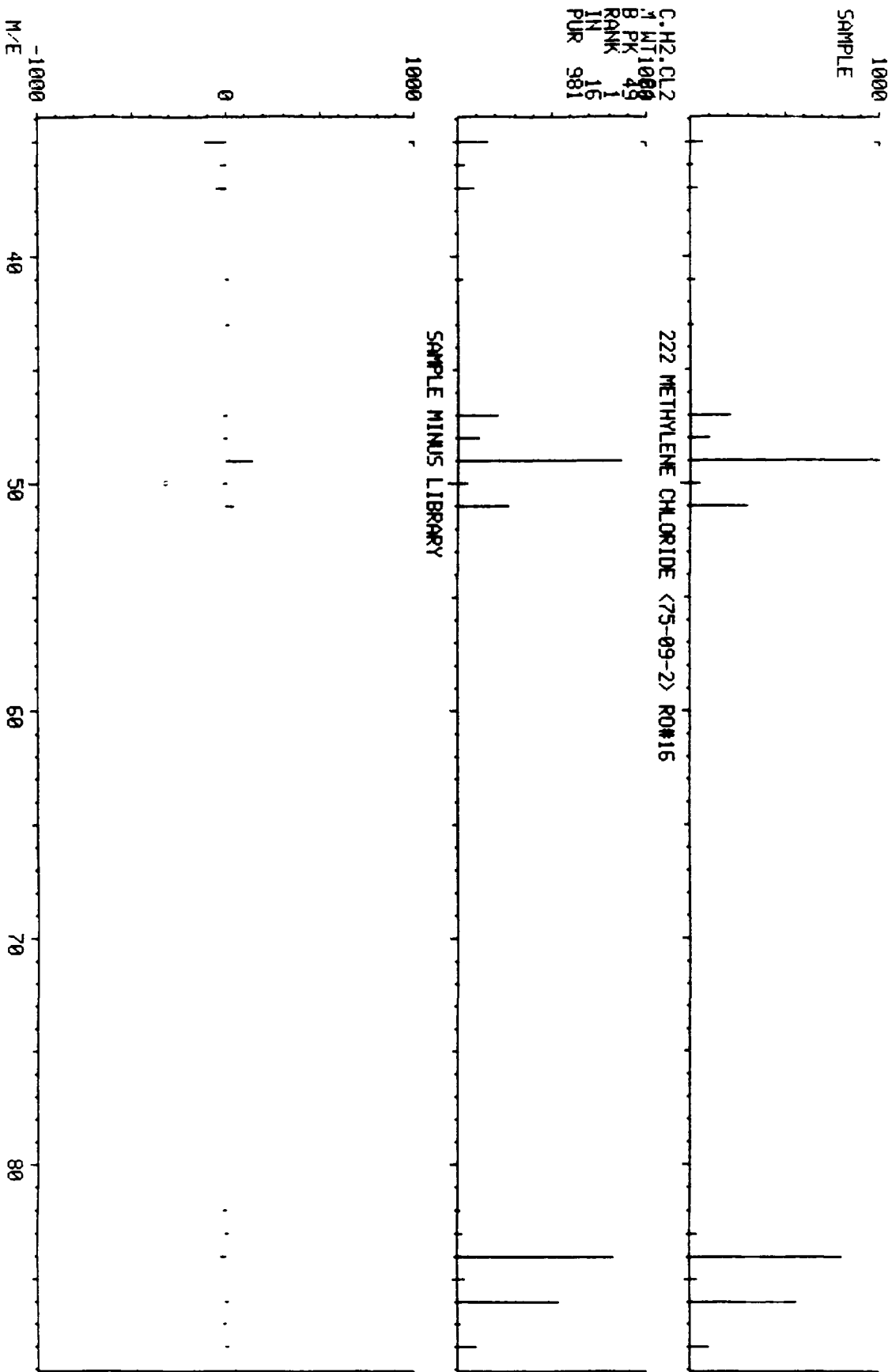
LIBRARY SEARCH
12/22/89 8:15:00 + 3:21
SAMPLE: 5 GRAMS CC#309690 EPA#B2020 CASE#18756.7 ON#19
ENHANCED (5 158 2N 0T)

COMPUJHEM LABS

DATA: GH009690A19 # 268

BASE M/E: 49
RIC: 17151.

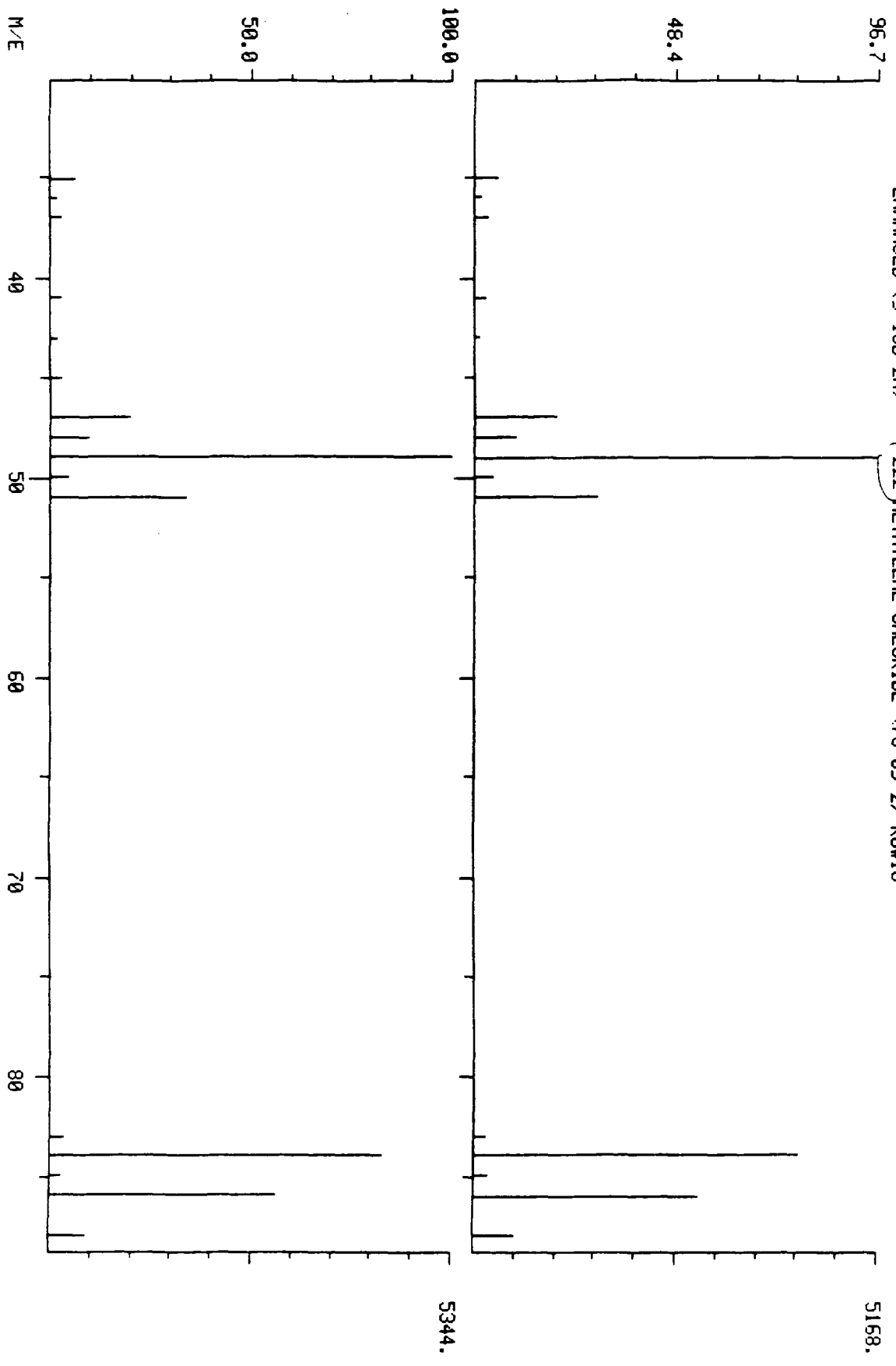
C-H2 CL2
I WT 1000
B PK 49
RANK 1
IN 16
PUR 981



DUAL MASS SPECTRUM
12/22/89 8:15:00 + 3:21
SAMPLE: 5 GRAMS CC#309690 EPA#: B202C CASE#18756.7 ON#19
ENHANCED (5 158 ZN) (222) METHYLENE CHLORIDE (75-09-2) RO#16

COMPUCHEM LABS

DATA: GH009690A19 #268 BASE M/E: 49/ 49
RIC: 17151./ 18143.



LIBRARY SEARCH
12/22/89 8:15:00 + 5:34
SAMPLE: 5 GRAMS CC#309690 EPA#: B202C CASE#18756.7 ON#19
ENHANCED (5 158 2N QT)

COMPUCHEM LABS

DATA: GH009690A19 # 446

BASE M/E: 43
RIC: 361.

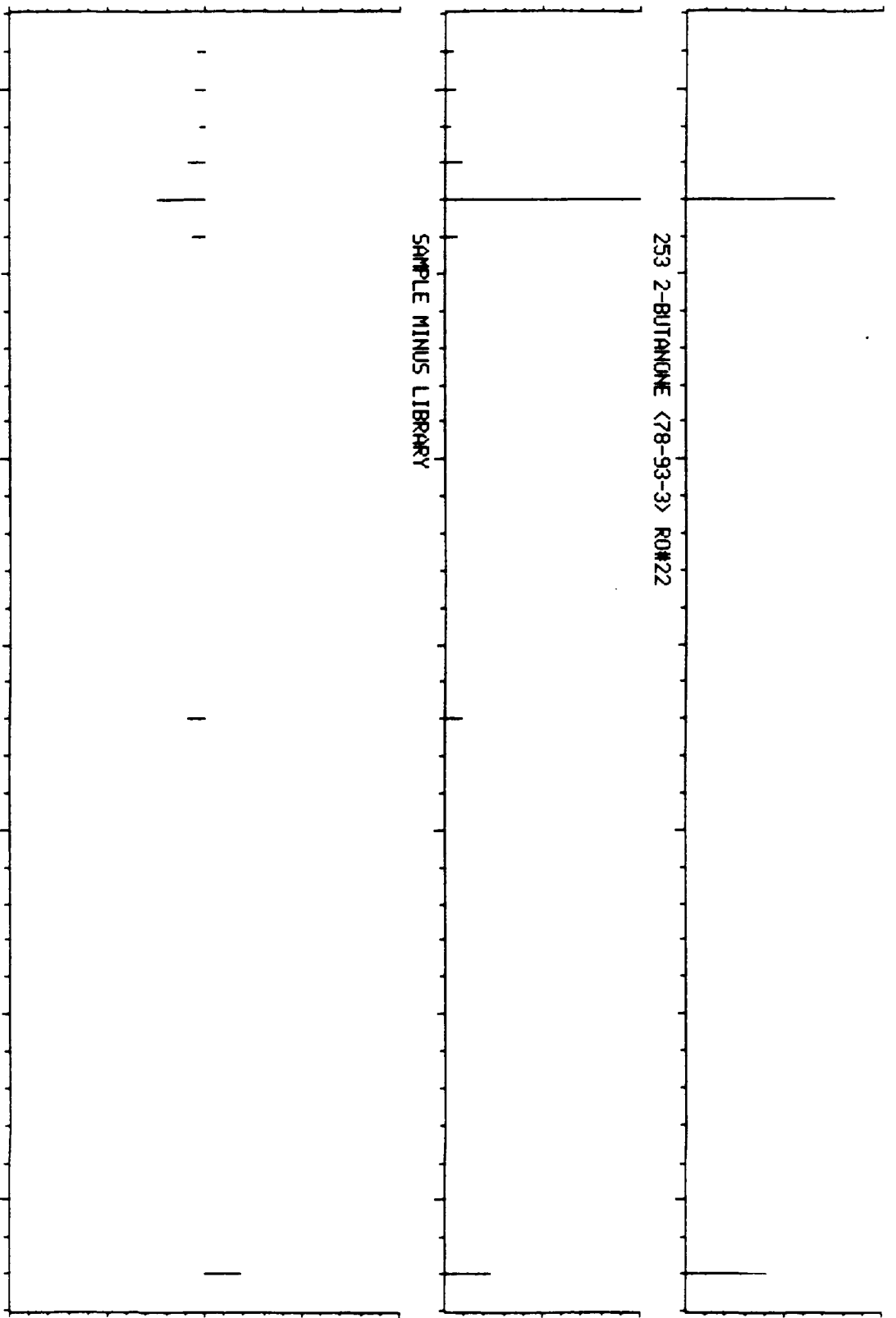
1322
SAMPLE

C4-H8-O
1 WT 1322
R PK 43
RANK 22
IN 1
PUR 764

253 2-BUTANONE <78-93-3> ROM#22

SAMPLE MINUS LIBRARY

-1322
M/E 40 45 50 55 60 65 70

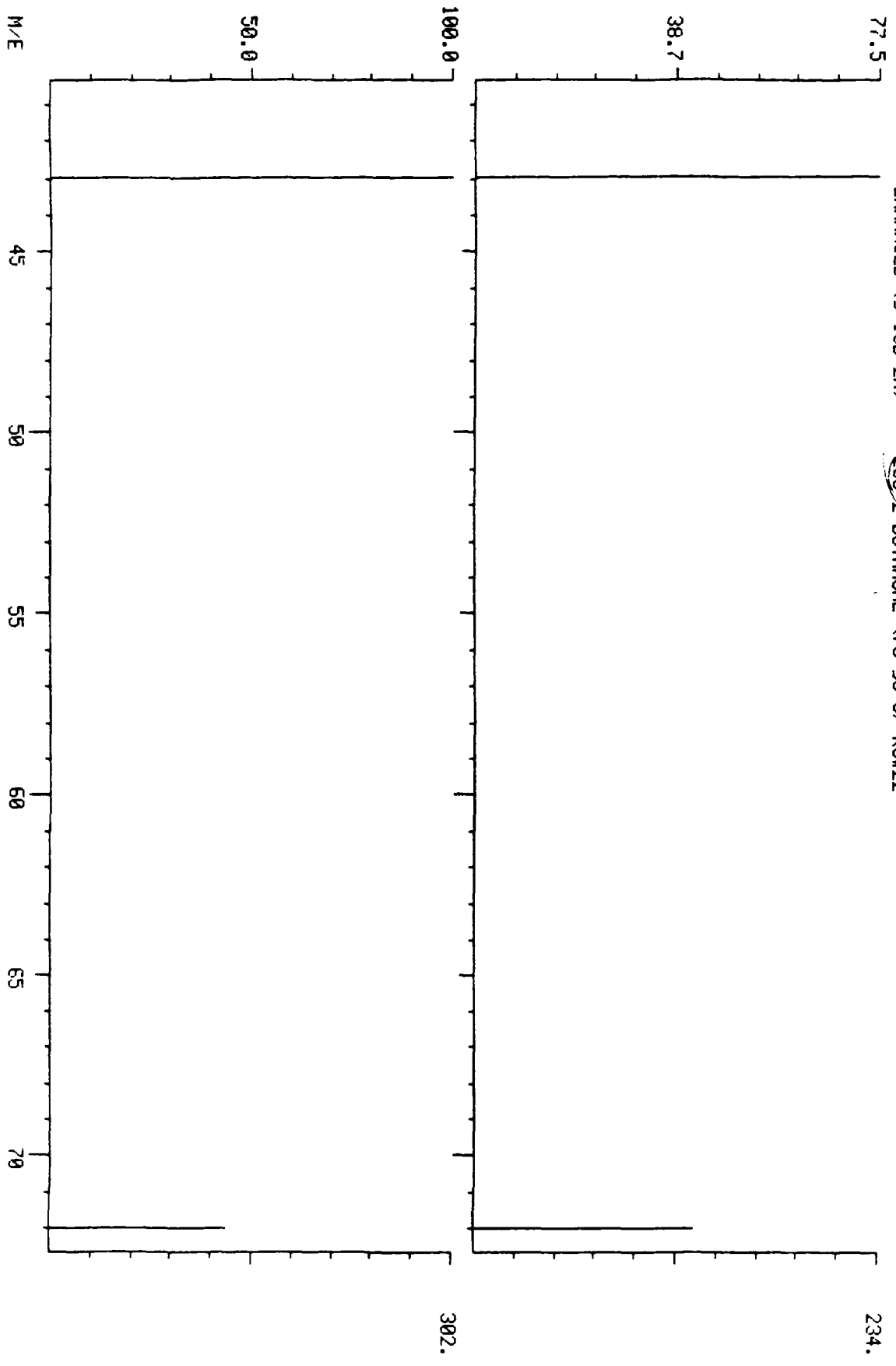


DUAL MASS SPECTRUM
12/22/89 8:15:00 + 5:34
SAMPLE: 5 GRAMS CC#309690 EPA#: B202C CASE#18756.7 ON#19
ENHANCED (5 158 2N) (53 2-BUTANONE (78-93-3) R0#22

COMPUCHEN LABS

DATA: GH009690A19 #446

BASE M/E: 43/ 43
RIC: 361. / 433.



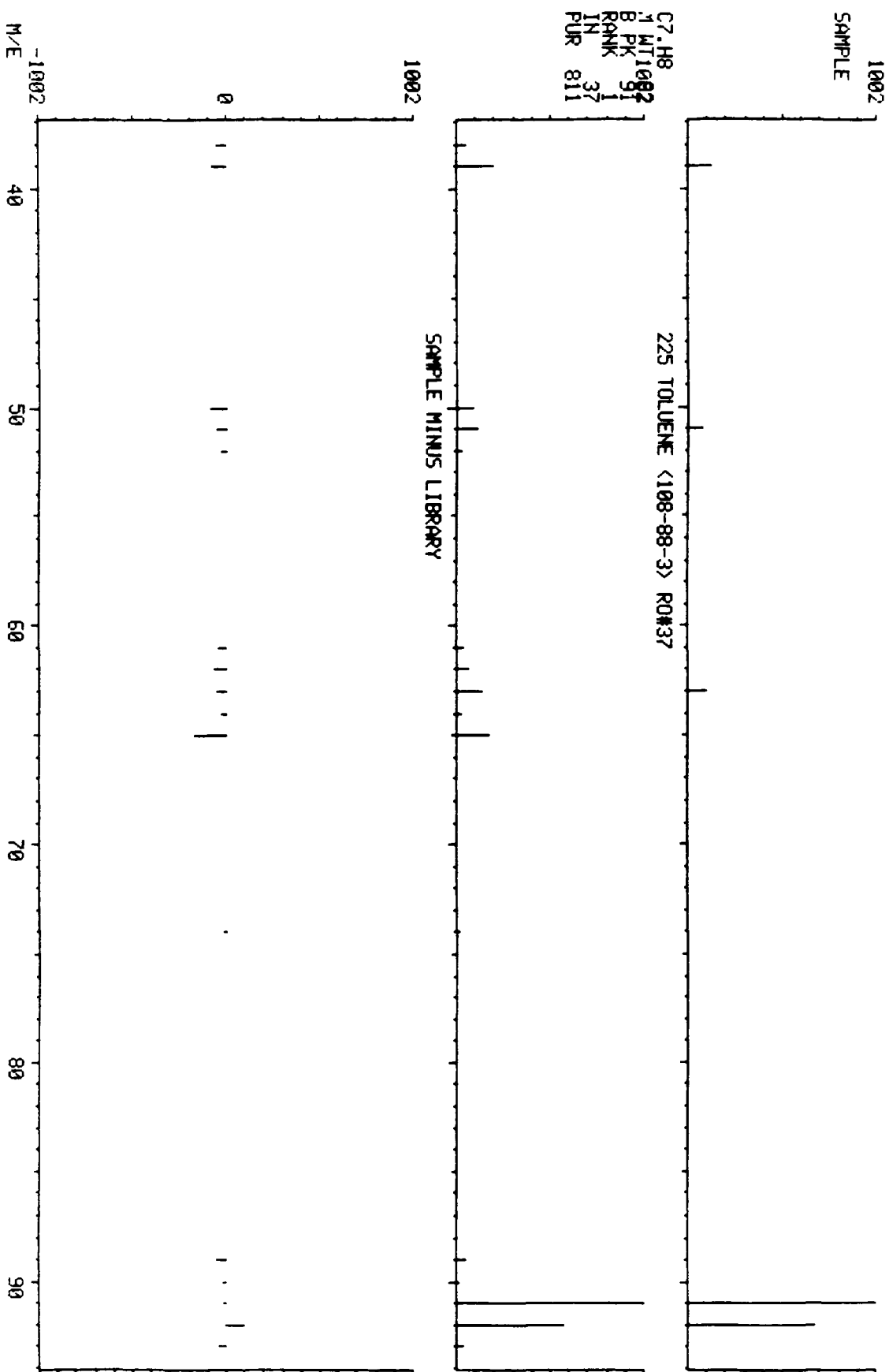
LIBRARY SEARCH
12/22/89 8:15:00 + 10:07
SAMPLE: 5 GRAMS CC#309690 EPA#: 62020 CASE#18756.7 ON#19
ENHANCED (5 158 2M 0T)

COMPUCHEM LABS

DATA: GH005690A19 # 809

BASE M/E: 91
RIC: 1277.

C7.H8
1 WT 1002
B PK 91
RANK 1
IN 37
PUR 811

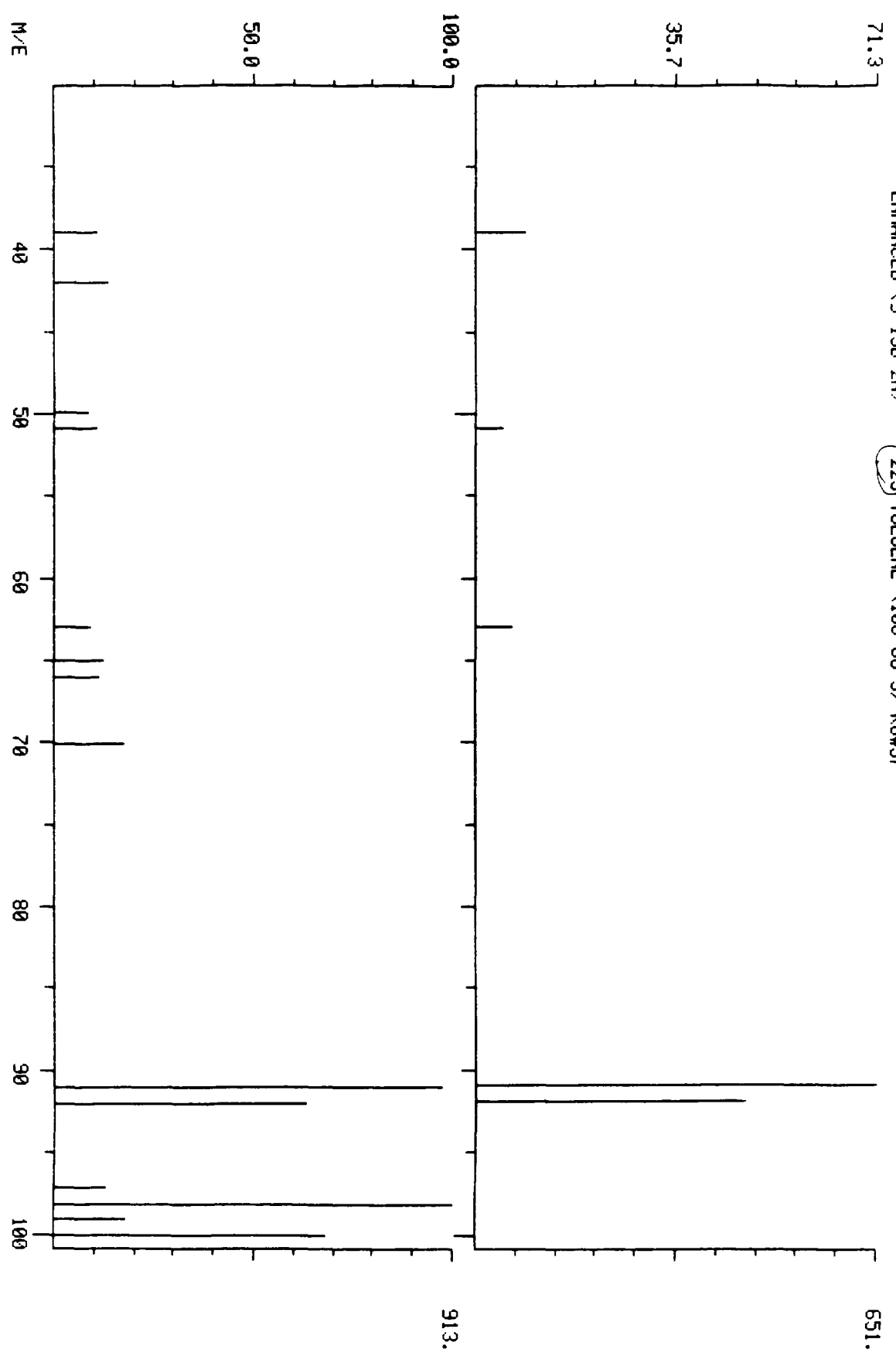


DUAL MASS SPECTRUM
12/22/89 8:15:00 + 10:07
SAMPLE: 5 GRAMS CC#309690 EPA#: B202C CASE#18756.7 ON#19
ENHANCED (5 158 2N) (225) TOLUENE (108-88-3) RM#37

COMPUCHEM LABS

DATA: GH009690A19 #809

BASE M/E: 91 / 98
RIC: 1277. / 4135.

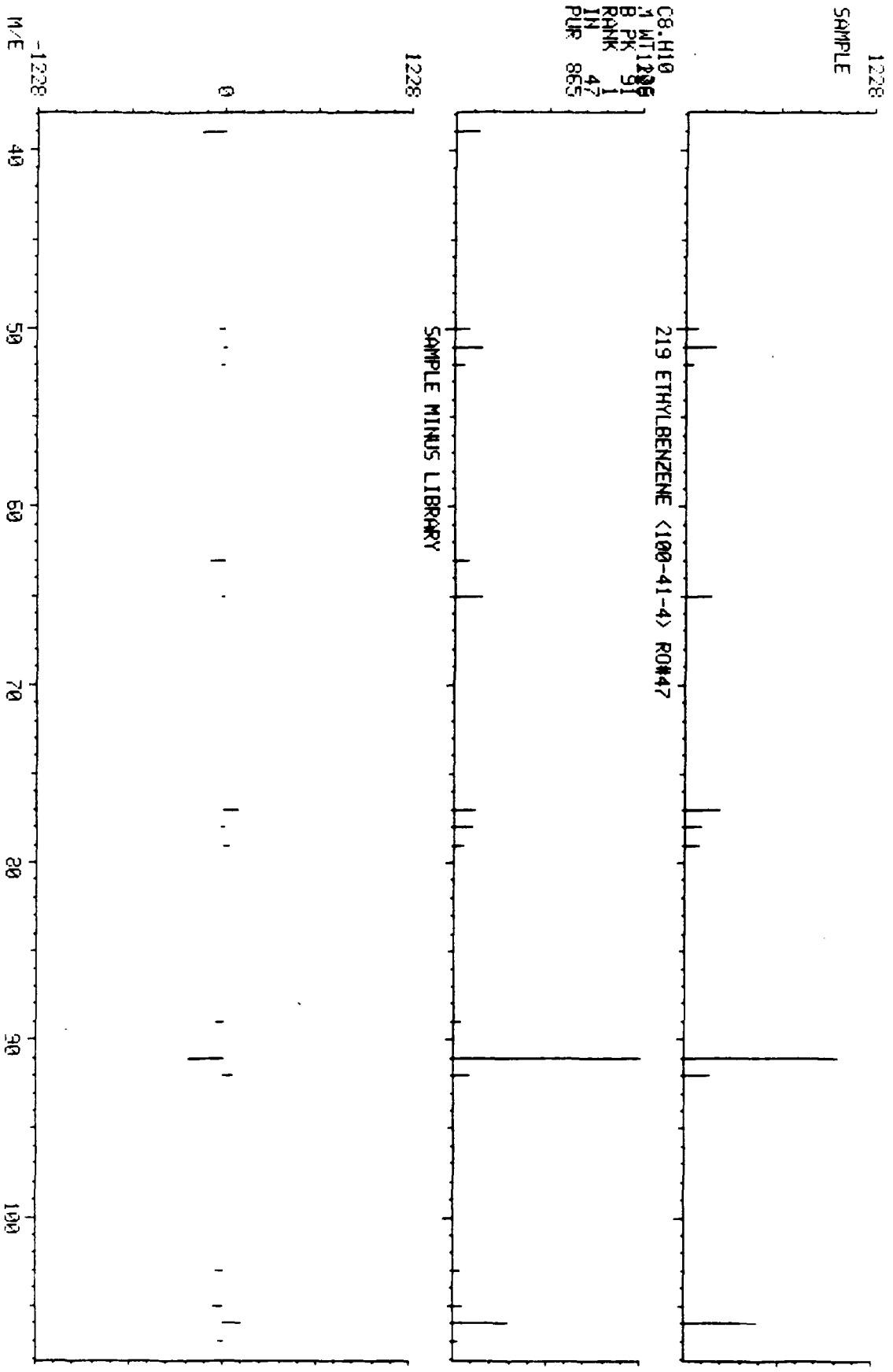


LIBRARY SEARCH
12/22/89 8:15:00 + 12:51
SAMPLE: 5 GRAMS CC#309690 EPA#:B2020 CASE#18756.7 QM#19
ENHANCED (5 158 2H 0T)

COMPUCHEM LABS

DATA: GH009690A19 #1028

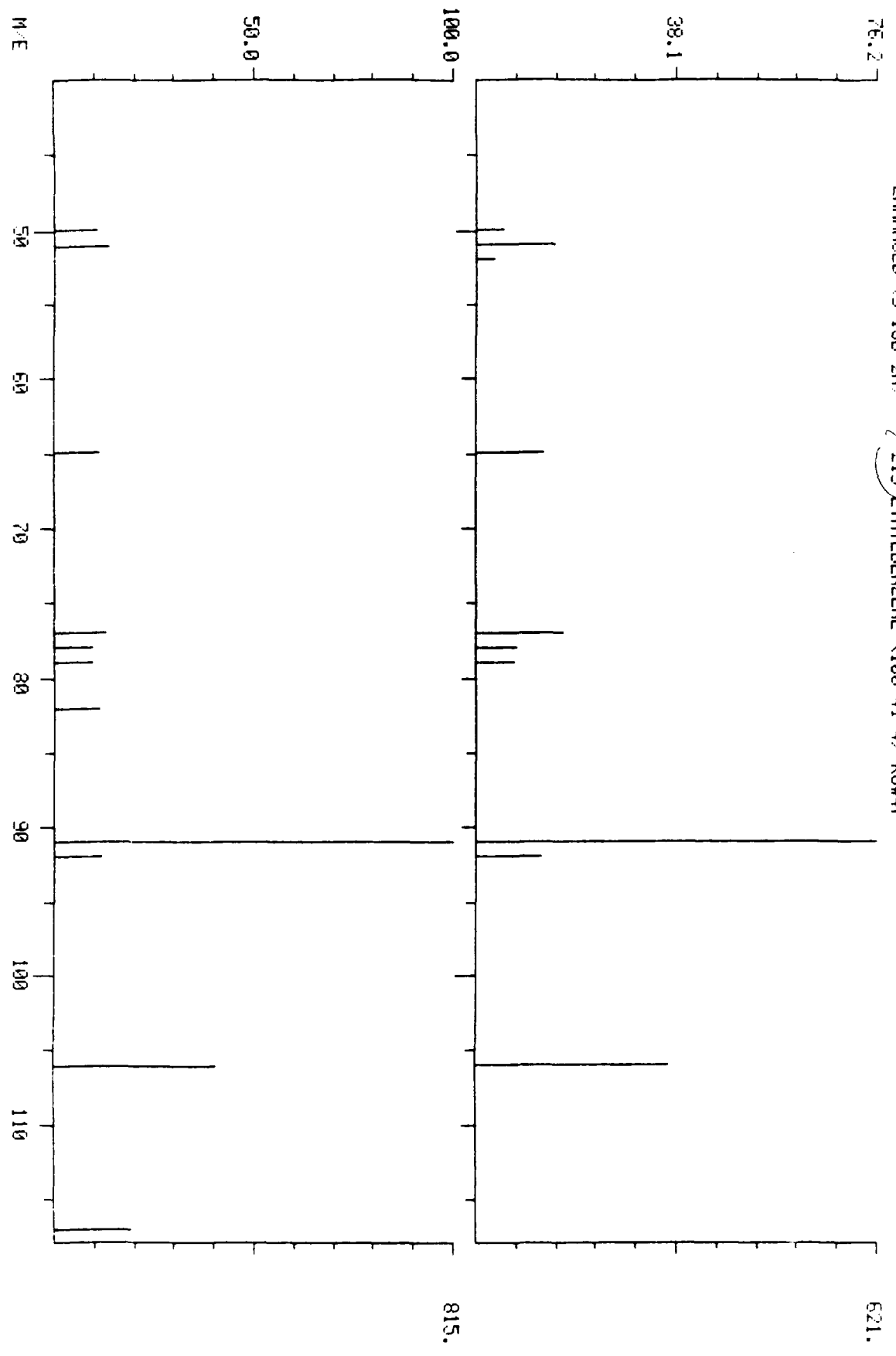
BASE M/E: 91
RIC: 1577.



DUAL MASS SPECTRUM
12/22/89 8:15:00 + 12:51
SAMPLE: 5 GRAMS CC#309690 EPA#: 82020 CASE#18756.7 QM#19
ENHANCED (5 158 ZN) 219 ETHYLBENZENE <100-41-4> R0#47

COMPUchem LABS

DATA: GH009690M19 #1028 BASE M/E: 91 / 91
R10: 1577. / 2035.

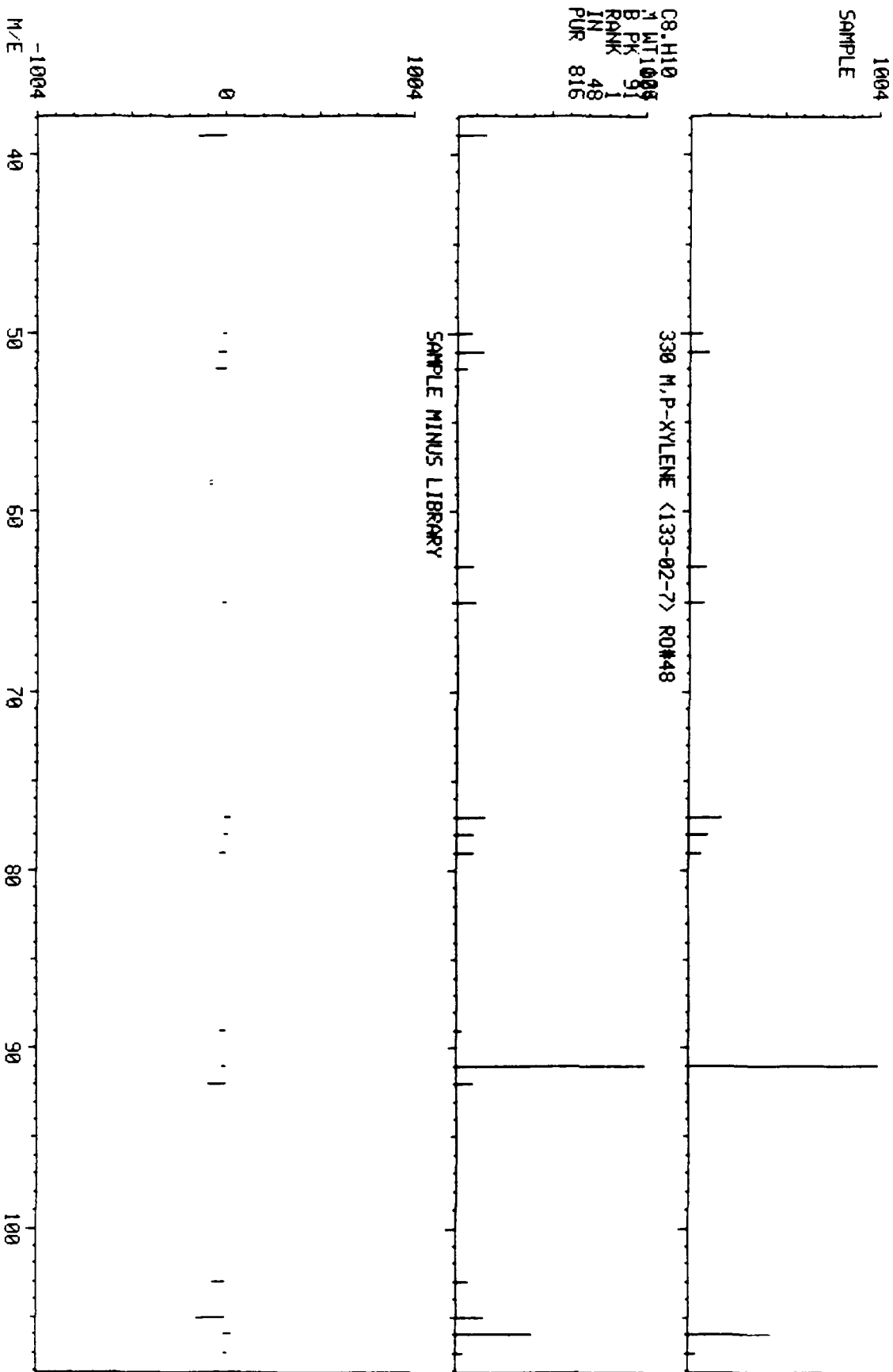


COMPUCHEM LABS
LIBRARY SEARCH
12/22/89 8:15:00 + 13:05
SAMPLE: 5 GRAMS CC#309690 EPA#: B202C CASE#18756.7 ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEM LABS

DATA: CH009690A19 #1048

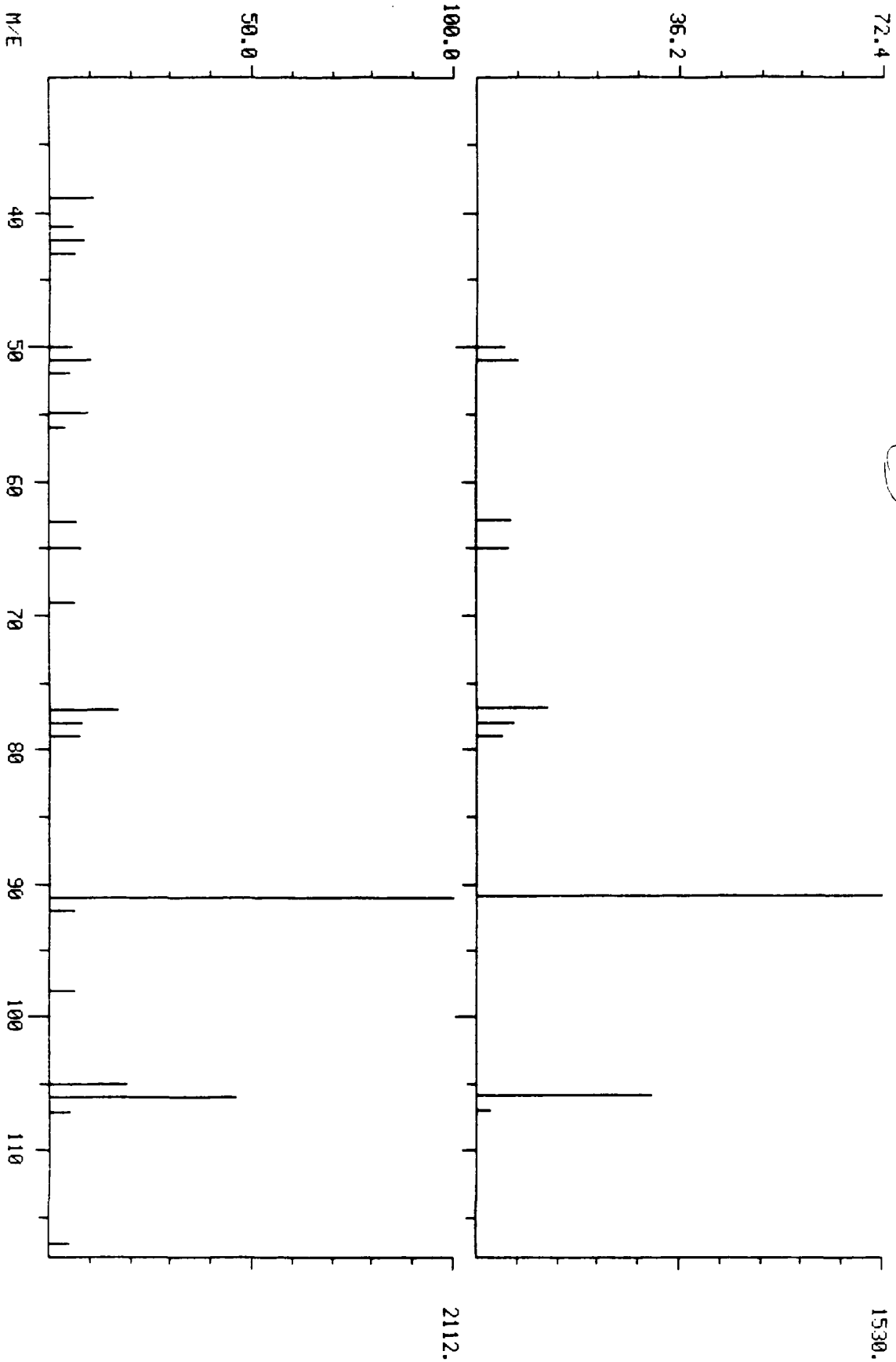
BASE M/E: 91
RIC: 3257.



DUAL MASS SPECTRUM
12/22/89 8:15:00 + 13:06
SAMPLE: 5 GRAMS CC#309690 EPA#: B202C CASE#18756.7 ON#19
ENHANCED (5.158 ZN) (330 M, P-XYLENE (133-02-7) R0#48

COMPUCHEM LABS

DATA: GH009690A19 #1048 BASE M/E: 91/ 91
RIC: 3267.7 6471.

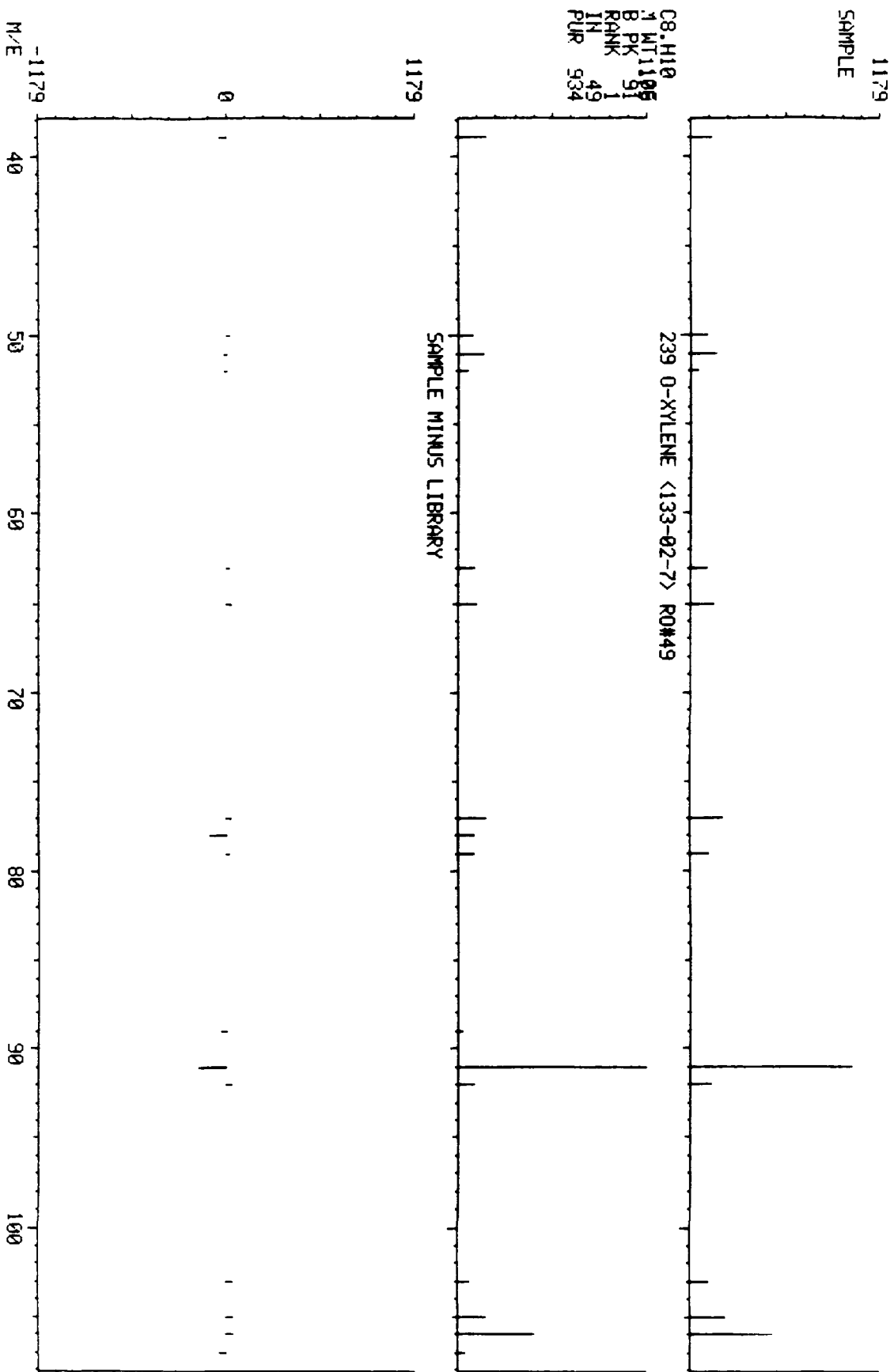


LIBRARY SEARCH
12/22/89 8:15:00 + 13:47
SAMPLE: 5 GRAMS CC#309690 EPA#:8202C CASE#18756.7 ON#19
ENHANCED (5.158 ZN 0T)

COMPUCHEM LAE5

DATA: GH009690A19 #1103

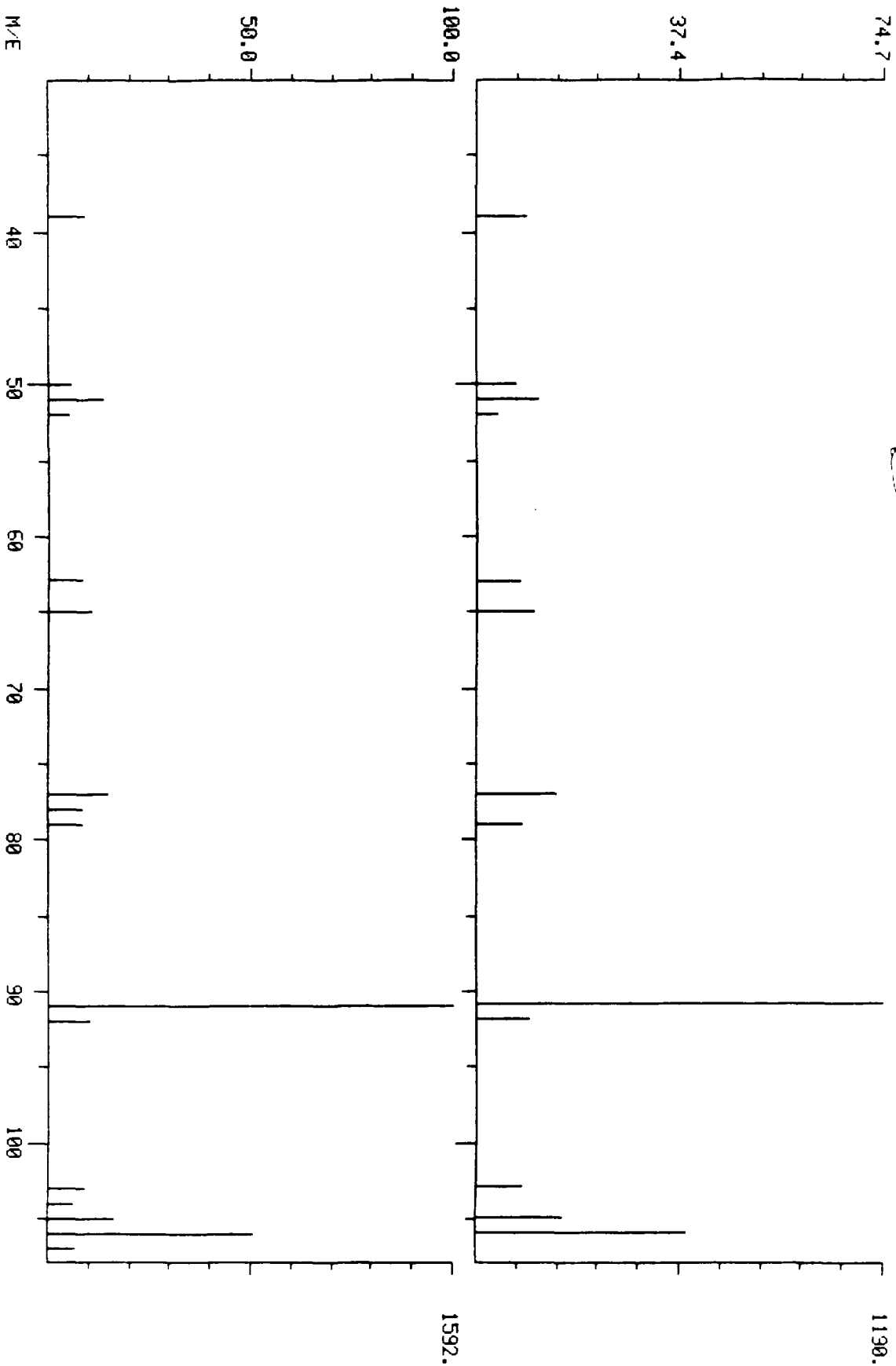
BASE M/E: 91
RIC: 3523.



DUAL MASS SPECTRUM
12/22/89 8:15:00 + 13:47
SAMPLE: 5 GRAMS CC#309690 EPA#: B202C CASE#18756.7 ON#19
ENHANCED (5 158 2N) (239)0-XYLENE (133-02-7) RM#49

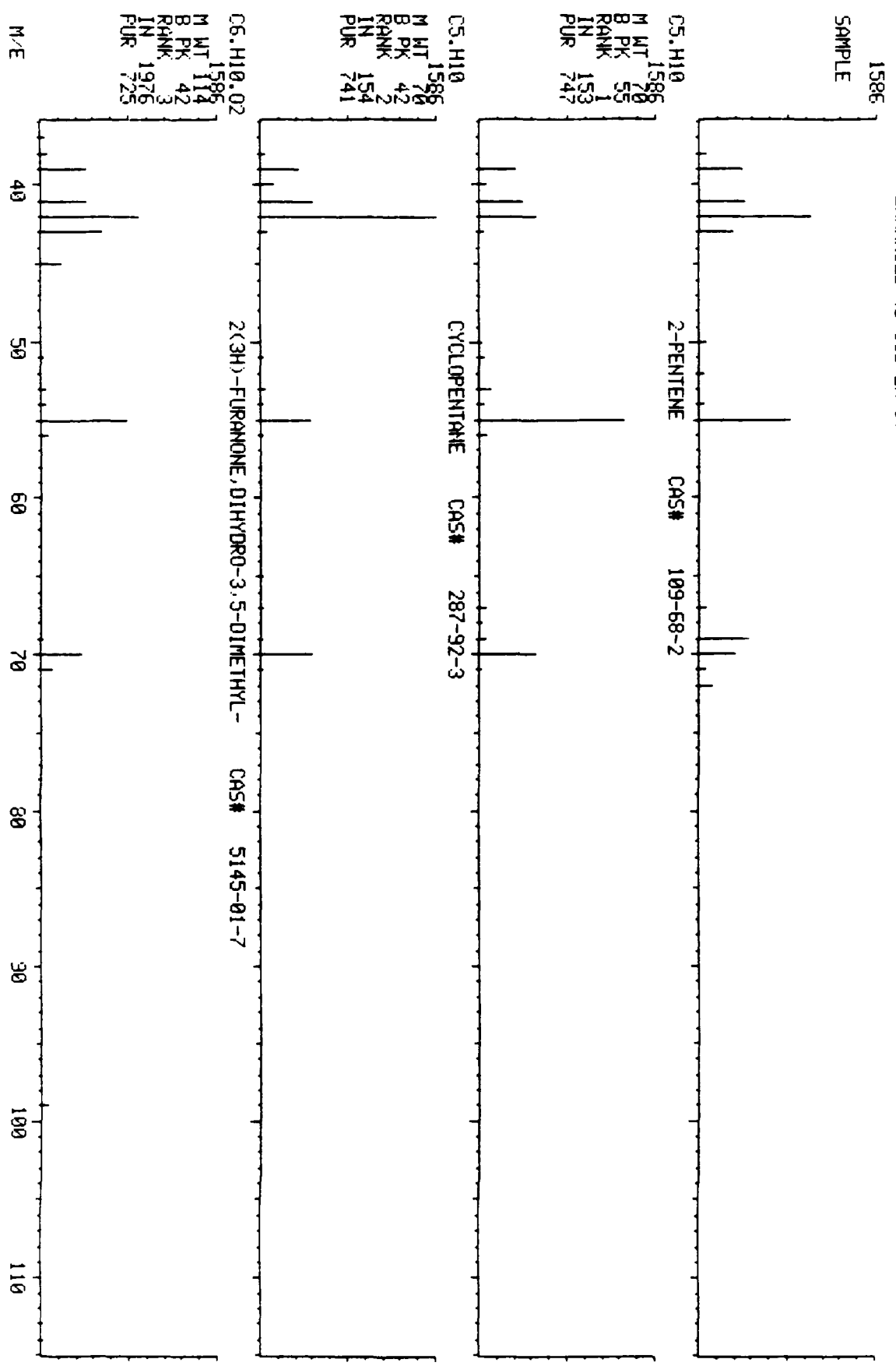
COMPUCHEM LABS

DATA: GH009690A19 #1103 BASE M/E: 91/ 91
RIC: 3523./ 4511.



LIBRARY SEARCH
 12/22/89 8:15:00 + 13:01
 SAMPLE: 5 GRAMS CC#309690 EPA#:B202C CASE#18756.7 ON#19
 ENHANCED (S 158 2N 0T)

COMPUCHEM LABS
 DATA: CH009690A19 #1042
 BASE M/E: 42
 RIC: 2899.



LIBRARY SEARCH
12/22/89 8:15:00 + 15:35
SAMPLE: 5 GRAMS CC#309690 EPA#: B2020 CASE#18756.7 ON#19
ENHANCED (5 158 2H 0T)

COMPUCHEM LABS

DATA: GH009690A19 #1247

BASE M/E: 84
RIC: 6671.

1000
SAMPLE

CYCLOPENTANONE, 2-ETHYL-

CAS# 4971-18-0

C7.H12.0
1000
M WT 112
3 PK 84
RANK 1
IN 1787
PUR 725

CYCLOHEPTANONE

CAS# 502-42-1

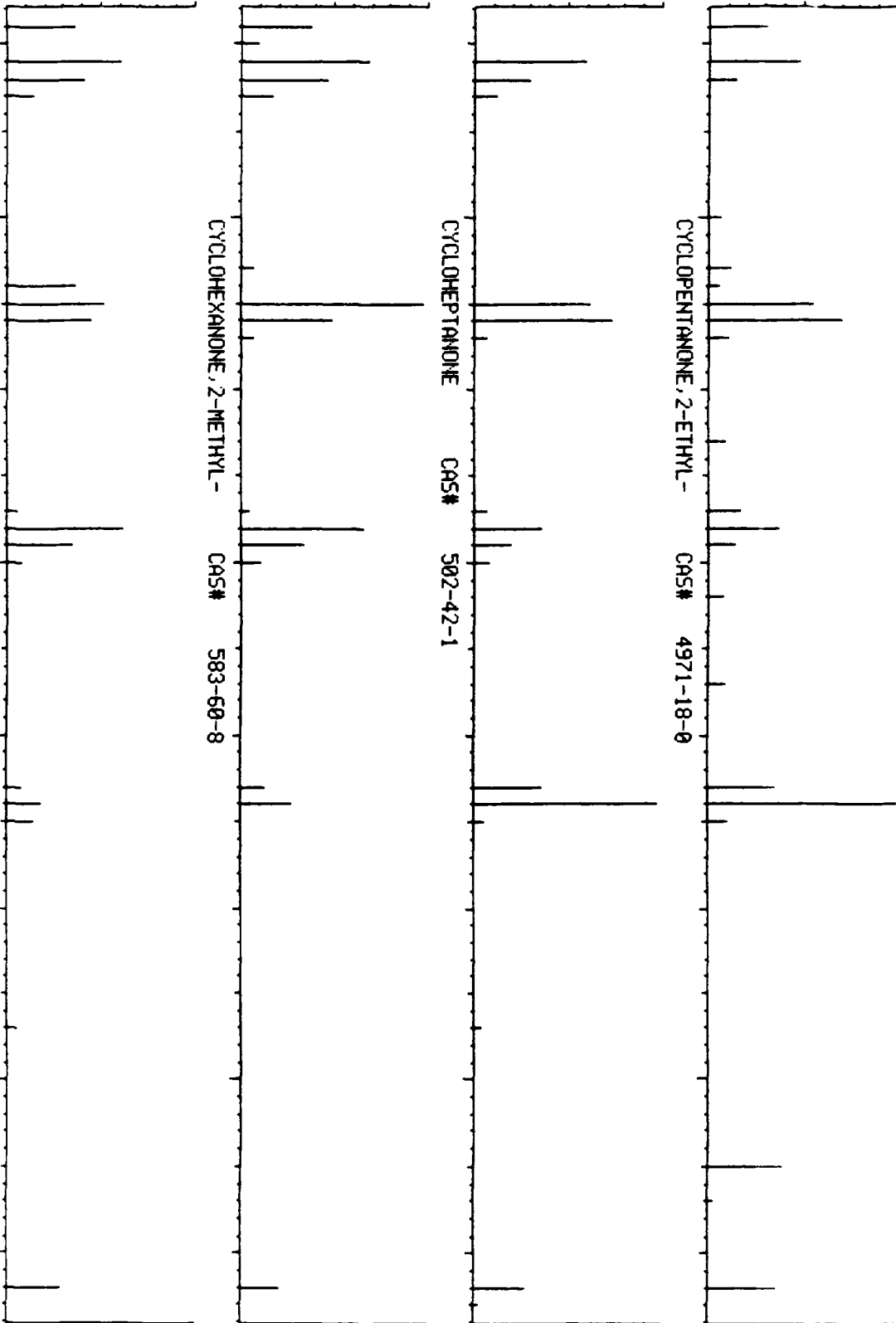
C7.H12.0
1000
M WT 112
5 PK 55
RANK 2
IN 1771
PUR 646

CYCLOHEXANONE, 2-METHYL-

CAS# 583-60-8

C7.H12.0
1000
M WT 112
6 PK 68
RANK 3
IN 1772
PUR 624

M/E 40 50 60 70 80 90 100 110



LIBRARY SEARCH
 12/22/89 8:15:00 + 15:58
 SAMPLE: 5 GRAMS CC#309690 EPA#: 8202C CASE#18756.7 ON#19
 ENHANCED (5 158 2H 0T)

COMPUCHEM LABS

DATA: GH009690A19 #1278

BASE M/E: 281
 RIC: 4967.

1006
 SAMPLE

C8.H24.04.514
 M WT 1006
 B PK 296
 RANK 281
 IN 21804
 PUR 828

CYCLOTETRASILOXANE, OCTAMETHYL - CAS# 556-67-2

C12.H36.04.515

M WT 1006
 B PK 384
 RANK 281
 IN 26593
 PUR 582

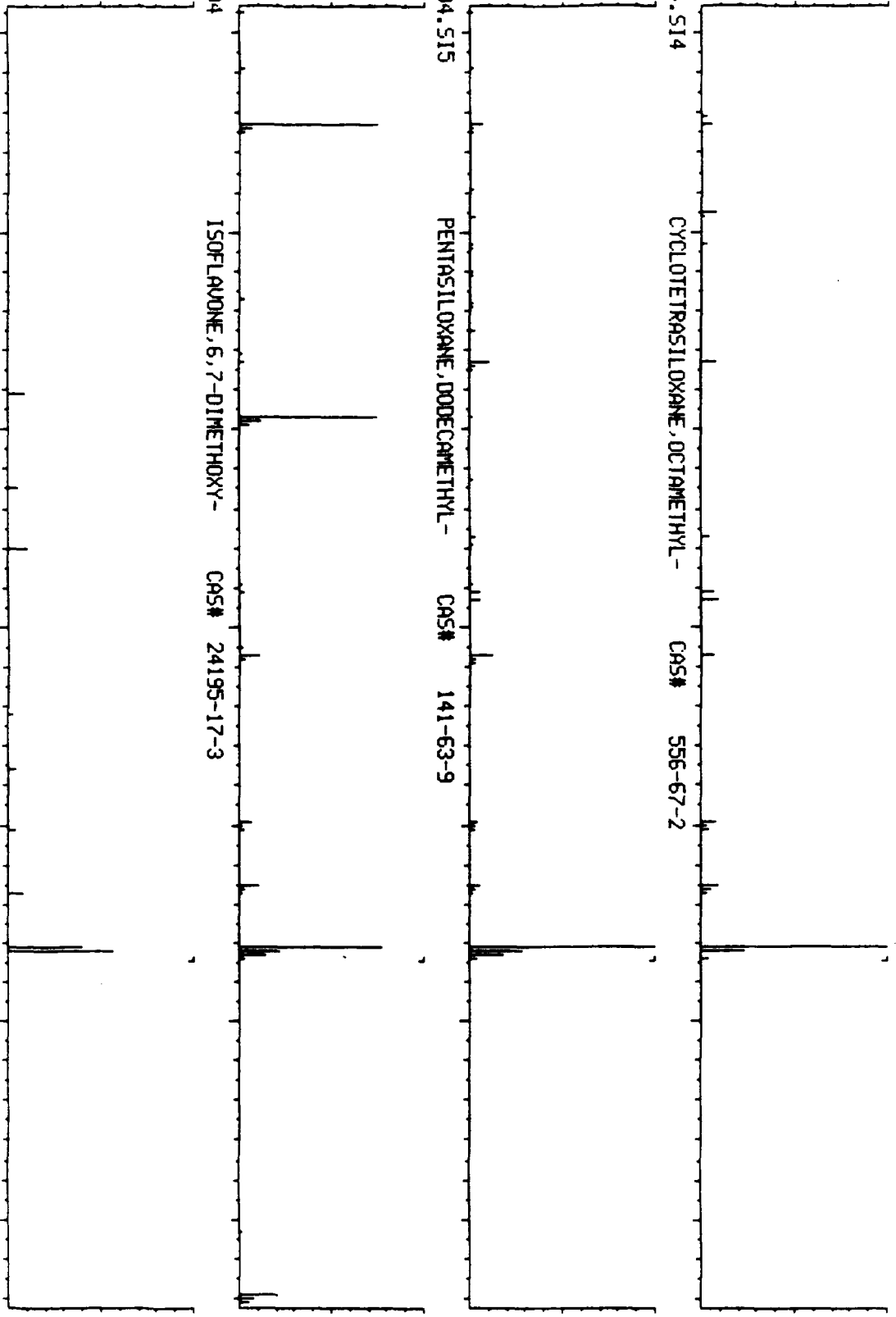
PENTASILOXANE, DODECAMETHYL - CAS# 141-63-9

C17.H14.04

M WT 1006
 B PK 282
 RANK 3
 IN 20667
 PUR 502

ISOFLAVONE, 6,7-DIMETHOXY - CAS# 24195-17-3

M/E 50 100 150 200 250 300 350



LAB INSTRUCTIONS:

RECEIPT DATE 12/20/89

CASE#: 18756 7

DUE DATE

VOA
GC/MS WORKSHEET

COMPUCHEM#: 309690

R1 [] R2 [] D1 [] ()
R3 [] R4 [] D2 [] ()

L L. SOLID, EPA SOW 2/88

Sample Prep Code---155
Instrument Code----413
Compound List-----494
Surrogate Std-----394
Internal Std-----036

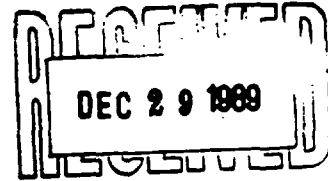
=====

SAMPLE ID# B202C Dry Wt. Factor 1.15 % Moisture 13

=====

GC/MS ANALYSIS

Amount Purged: [] 10mls/Xg soil or [] Dilution _____ ul/10000ul/Xg soil
Internal Standard Volume Added 5 ul
Surrogate Standard Volume Added 5 ul
BFB Filename B6891221B19 Disk ()
Blank Filename G4009785C19 Disk ()
Standard Filename G3891221B19 Disk ()
Sample Filename G4009690A19 Disk ()



ANALYST(S): Injection M92/Alan Frank

Work-up 1452

GC/MS REVIEW

CONDITION
CODE

OK

Entry Codes OK, EA, ES, SM, JS, SL, SH, JA, DA

Non-Entry Codes IM, IL, IH, SW, CT, CS, PC, NR
IF, LA, DI, CO, RN, DW, SI, SF
UP, BB, OT, VC, FO, NS

Disposition: [] Complete
[] Reprep neat required
[] Reprep using _____ g
[] Dilute (:)

Extraneous Peak Search Results:
of Peaks Found: 03

Quality Assurance Notice(s):
Notices Required 1



COMMENTS:

GC/MS Review [Signature] Date 12/29/89 Auditor DWagner Date 12/29/89

REPORT INTEGRATION

Total # of Injections 1

Final Reportable Package(s): GHO-A19

QA COMMENTS:

=====

INITIALS _____ DATE _____/_____/_____
=====

FINAL REVIEW INITIALS _____ DATE _____/_____/_____

AC1007 (05/89)

VOLATILE PREPARATION WORKSHEET

QUEUE # 19 PREP CODE -155 ASSIGNED TO Harshad Jeshi DATE 12/20/89

SAMPLE NUMBER	CASE NUMBER	SDG	QC SAMPLE		SAMPLE WEIGHT (G) VOLUME (ML)	SAMPLE ID	COMMENTS
			TYPE	ORIGINAL			
309763	18520	0164			5.0g	TB-12	
309768	↓	↓			5.0g	WB5-1	
309769	↓	↓			5.0g	WB5-2	
309770	↓	↓			5.0g	WB5-4	
309679	18756	0007			5.0g	B201A	
309680	↓	↓	SS	309679	5.0g	B201A MS	
309681	↓	↓	SS	309679	5.0g	B201A MSD	
309682	↓	↓	BS		0.0g		
309686	↓	↓			5.0g	B201B	
309687	↓	↓			5.0g	B202AR	
309688	↓	↓			5.0g	B202A	
309689	↓	↓			5.0g	B202B	
309690	↓	↓			5.0g	B202C	
309784			B1		5.0 ml	B3	
309785			B2		0.0 ml	B4	
309786			B3		0.0 ml	B5	
309787			B4		0.0 ml	B6	
309788			B5		0.0 ml	B7	

SURROGATE # LOT # MANUAL OPERATOR 730, 516
 AMOUNT
 RELINQUISHED BY [Signature] DATE 12/20/89 RECEIVED BY [Signature] DATE 12/20/89

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	462	49100	50.0		X 115
221	50	CHLOROMETHANE				BDL	24
231	62	VINYL CHLORIDE				BDL	1
220	94	BROMOMETHANE				BDL	1
209	64	CHLOROETHANE				BDL	1
216	96	1,1-DICHLOROETHENE				BDL	1
254	76	CARBON DISULFIDE				BDL	1
252	43	ACETONE (2-PROPANONE)			22.8	23 20.8	1
248	114 I	1,4-DIFLUOROBENZENE (IS)	608	202000	50.0		
222	84	METHYLENE CHLORIDE			19.6	20 23.8	1
226	96	TRANS-1,2-DICHLOROETHENE				BDL	1
214	63	1,1-DICHLOROETHANE				BDL	1
257	43	VINYL ACETATE				BDL	1
237	96	CIS-1,2-DICHLOROETHENE				BDL	1
253	72	2-BUTANONE			3.0	34 35	1
211	83	CHLOROFORM				BDL	1
227	97	1,1,1-TRICHLOROETHANE				BDL	1
206	117	CARBON TETRACHLORIDE				BDL	1
203	78	BENZENE				BDL	1
215	62	1,2-DICHLOROETHANE				BDL	1
270	117 I	D5-CHLOROETHANE (IS) RO#29	999	211000	50.0		
229	130	TRICHLOROETHENE				BDL	1
217	63	1,2-DICHLOROPROPANE				BDL	1
212	83	BROMODICHLOROMETHANE				BDL	1
218	75	CIS-1,3-DICHLOROPROPENE				BDL	1
256	43	4-METHYL-2-PENTANONE				BDL	1
225	92	TOLUENE			1.7	20 25	1
250	75	TRANS-1,3-DICHLOROPROPENE				BDL	1
228	97	1,1,2-TRICHLOROETHANE				BDL	1
224	164	TETRACHLOROETHENE				BDL	1
255	43	2-HEXANONE			3.2	BDL 30 45	10
208	129	DIBROMOCHLOROMETHANE				BDL	1
207	112	CHLOROBENZENE				BDL	1
219	106	ETHYLBENZENE			1.2 4.4	45 45	1
330	106	M, P-XYLENE			3.2 32	30 45	1
239	106	O-XYLENE			2.8	30 35	1
251	104	STYRENE				BDL	1
205	173	BROMOFORM				BDL	1
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	1
258	65 S	D4-1,2-DICHLOROETHANE RO#57			48.3	97. %	1
247	95 S	BROMOFLUOROBENZENE			45.4	91. %	1
233	98 S	D8-TOLUENE RO#59			47.0	94. %	1
289	106	XYLENES (TOTAL)			6.0	67	1

CORRECTED/REVIEWED BY C. J. Anderson
(GC/MS DATA REVIEWER)

DATE 12-27-85

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
299	96	1,2-DICHLOROETHENE (TOTAL)				BDL	
CHECKSUMS:							
		3979.	2069	462100.		357.4	349.

CORRECTED/REVIEWED BY AKL
(GC/MS DATA REVIEWER)

DATE 12 27 85

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P
40	258	D4-1,2-DICHLOROETHANE RO#57	48.3	50.0	97.	70-121	X
41	247	BROMOFLUOROBENZENE	45.4	50.0	91.	74-121	X
42	233	DB-TOLUENE RO#59	47.0	50.0	94.	81-117	X

* ADVISORY SURROGATE ONLY

++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

=====

CORRECTION FACTOR CALCULATION:

$$\frac{5.0 \text{ G}}{\text{WET WEIGHT OF SAMPLE (G)}} \times \frac{\text{GC/MS DILUTION FACTOR}}{\text{GC/MS DILUTION FACTOR}} \times \text{DRY WEIGHT FACTOR} =$$

$$\frac{5.0 \text{ G}}{5.00 \text{ (G)}} \times \frac{1.0}{1.0} \times \frac{1.15}{1.00} = \frac{1.15}{1.00}$$

=====

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:

THE SURROGATES ARE ADDED TO THE SAMPLE PRIOR TO SPARGING.
SURROGATE SPIKE CONVERSION FACTOR = 1.

=====

VERSION 8

CORRECTED/REVIEWED BY C. E. St. Hill
(GC/MS DATA REVIEWER)

DATE 12-22-89

LABORATORY NOTICE

Sample I.D. # 309690

Client I.D. # B202C

CASE # 18756.7

The analysis of the volatile fraction of this sample indicated the presence of one or more extraneous peaks at the scan(s) listed below which were tentatively identified as siloxanes. This is a frequently observed laboratory artifact and should not be considered an actual sample constituent. It is being reported as an instrument artifact. While the peak(s) have been included in the total number of Library Searches, they have not been counted as one of the ten most intense TIC's requiring Library Search. If the siloxane was also detected in the associated blank, it has been flagged in the sample data with a "B", whether or not its concentration in the blank was great enough to require a Library Search. The data is being reported with reference to this qualifier.

scan(s): 1278 _____

Data Reviewer: *C. H. Stett*

Date: 12-27-89

Approved by: Robert J. Whitehead
Manager, Quality Assurance

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TAR

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Matrix: (soil/water) SOIL Lab Sample ID: 309687

Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009687C19

Level: (low/med) LOW Date Received: 12/20/89

% Moisture: not dec. 64 Date Analyzed: 12/27/89

Column: (pack/cap) CAP Dilution Factor: 1.0

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	140	U
74-83-9	-----Bromomethane	140	U
75-01-4	-----Vinyl Chloride	140	U
75-00-3	-----Chloroethane	140	U
75-09-2	-----Methylene Chloride	8300	EB
67-64-1	-----Acetone	3500	EB
75-15-0	-----Carbon Disulfide	69	U
75-35-4	-----1,1-Dichloroethene	69	U
75-34-3	-----1,1-Dichloroethane	69	U
540-59-0	-----1,2-Dichloroethene (total)	69	U
67-66-3	-----Chloroform	69	U
107-06-2	-----1,2-Dichloroethane	69	U
78-93-3	-----2-Butanone	5000	E
71-55-6	-----1,1,1-Trichloroethane	69	U
56-23-5	-----Carbon Tetrachloride	69	U
108-05-4	-----Vinyl Acetate	140	U
75-27-4	-----Bromodichloromethane	69	U
78-87-5	-----1,2-Dichloropropane	69	U
10061-01-5	-----cis-1,3-Dichloropropene	69	U
79-01-6	-----Trichloroethene	69	U
124-48-1	-----Dibromochloromethane	69	U
79-00-5	-----1,1,2-Trichloroethane	69	U
71-43-2	-----Benzene	1200	U
10061-02-6	-----Trans-1,3-Dichloropropene	69	U
75-25-2	-----Bromoform	69	U
108-10-1	-----4-Methyl-2-Pentanone	1200	U
591-78-6	-----2-Hexanone	12000	EB
127-18-4	-----Tetrachloroethene	69	U
79-34-5	-----1,1,2,2-Tetrachloroethane	69	U
108-88-3	-----Toluene	8600	E
108-90-7	-----Chlorobenzene	69	U
100-41-4	-----Ethylbenzene	8000	E
100-42-5	-----Styrene	2300	U
1330-20-7	-----Total Xylenes	15000	E

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

B202TAR

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 1.0 (g/mL) G Lab File ID: GR009687C19
 Level: (low/med) LOW Date Received: 12/20/89
 % Moisture: not dec. 64 Date Analyzed: 12/27/89
 Column (pack/cap) CAP Dilution Factor: 1.0

Number TICs found: 10

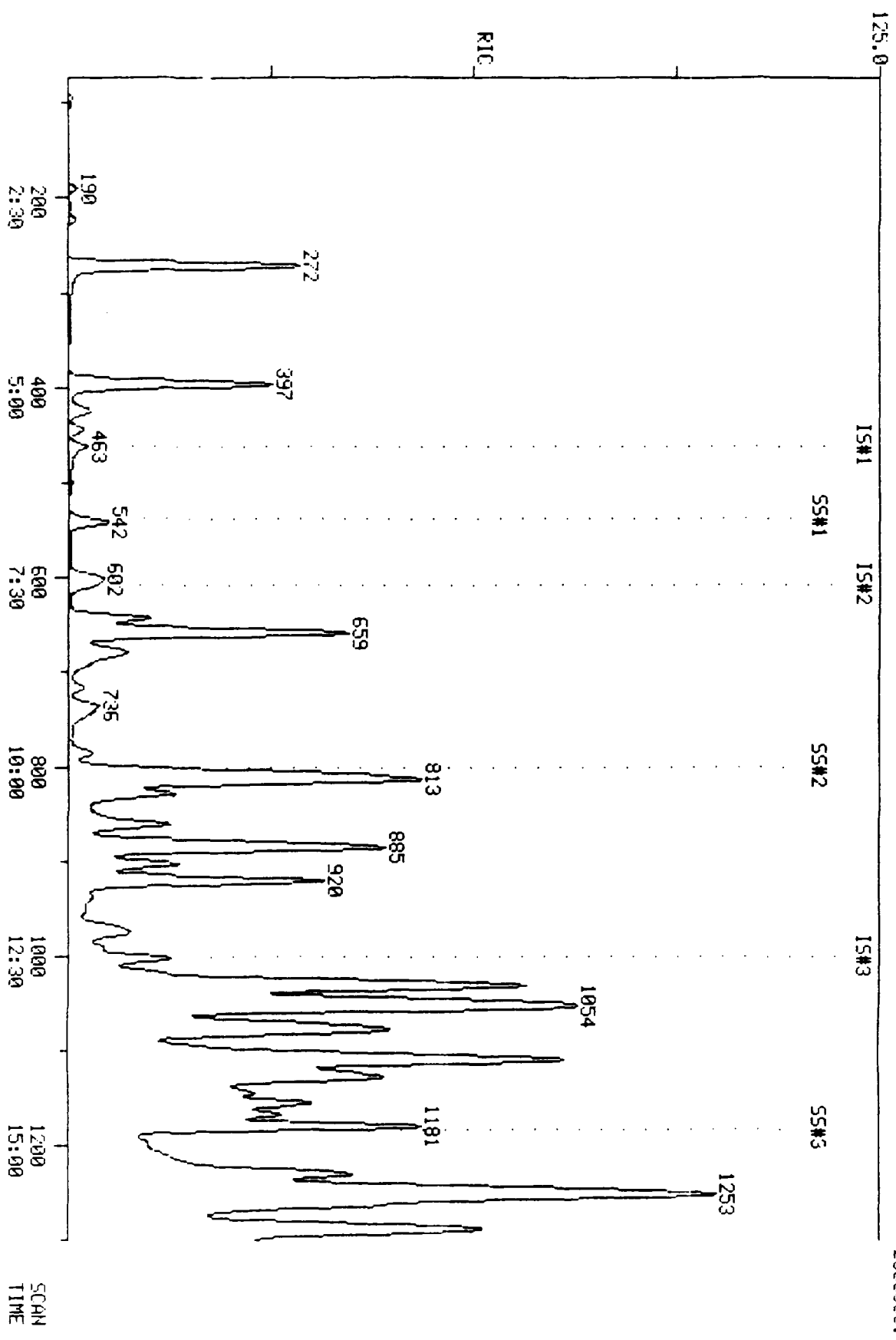
CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 534-22-5	FURAN, 2-METHYL-	4.95	9000	J
2. 625-86-5	FURAN, 2, 5-DIMETHYL-	8.23	13000	J
3.	UNKNOWN	11.05	4700	J
4.	UNKNOWN	11.50	3200	J
5.	UNKNOWN	13.49	2600	J
6.	UNKNOWN	14.09	2900	J
7. 13368-65-5	CYCLOHEXANONE, 3-METHYL-, (R)-	14.45	1700	J
8.	UNKNOWN	14.75	3900	J
9.	UNKNOWN	15.65	11000	J
10. 3141-02-4	1, 3-CYCLOPENTADIENE, 5-(1-MET	16.10	4400	J

RIC
12/27/89 7:10:00
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR OH#19
COND:.

COMPUchem LABS
COMPUchem DATA: GR009687C19 SCANS 75 TO 1300

2608630.



QUANTITATION REPORT FILE: GRC09687C19
 DATA: GRC09687C19.TI
 12/27/89 7:10:00
 SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR DN#19
 CONDS.:
 SUBMITTED BY: 19 ANALYST: 1422

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
 RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1
2	221 CHLOROMETHANE <74-87-3> RO#2
3	231 VINYL CHLORIDE <75-01-4> RO#3
4	220 BROMOMETHANE <78-83-9> RO#4
5	209 CHLOROETHANE <75-00-3> RO#5
6	216 1,1-DICHLOROETHENE <75-35-4> RO#8
7	254 CARBON DISULFIDE <75-15-0> RO#9
8	252 ACETONE (2-PROPANONE) <67-64-1> RO#13
9	*248 1,4-DIFLUOROBENZENE (IS) <540-36-3> RO#14
10	222 METHYLENE CHLORIDE <75-09-2> RO#16
11	226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17
12	214 1,1-DICHLOROETHANE <75-34-3> RO#19
13	257 VINYL ACETATE <108-05-4> RO#20
14	237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21
15	253 2-BUTANONE <78-93-3> RO#22
16	211 CHLOROFORM <67-66-2> RO#23
17	227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24
18	206 CARBON TETRACHLORIDE <56-23-5> RO#25
19	203 BENZENE <71-43-2> RO#26
20	215 1,2-DICHLOROETHANE <107-06-2> RO#27
21	*270 D5-CHLOROBENZENE (IS) RO#29
22	229 TRICHLOROETHENE <79-01-6> RO#30
23	217 1,2-DICHLOROPROPANE <78-87-5> RO#31
24	212 BROMODICHLOROMETHANE <75-27-4> RO#33
25	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35
26	256 4-METHYL-2-PENTANONE <108-01-1> RO#36
27	225 TOLUENE <108-88-3> RO#37
28	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38
29	228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39
30	224 TETRACHLOROETHENE <127-18-4> RO#41
31	255 2-HEXANONE <591-78-6> RO#42
32	208 DIBROMOCHLOROMETHANE <124-48-1> RO#43
33	207 CHLOROBENZENE <108-90-7> RO#45
34	219 ETHYLBENZENE <100-41-4> RO#47
35	330 M,P-XYLENE <133-02-7> RO#48
36	239 O-XYLENE <133-02-7> RO#49
37	251 STYRENE <100-42-5> RO#50
38	205 BROMOFORM <75-25-2> RO#51
39	223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54
40	*258 D4-1,2-DICHLOROETHANE RO#57
41	*247 BROMOFLUOROBENZENE <460-00-4> RO#58
42	*233 D8-TOLUENE RO#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	463	5:47	1	1.000	A BB	68062.	50.000 UG/KG	1.00
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	223	2:47	1	0.482	A BB	106545.	249.872 UG/KG	5.00
9	114	608	7:36	9	1.000	A BB	244667.	50.000 UG/KG	1.00
10	84	272	3:24	1	0.587	A BB	1484160.	599.342 UG/KG	11.98
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	445	5:34	1	0.961	A BB	66327.	362.147 UG/KG	7.24
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	543	6:47	9	0.893	A BB	429611.	88.102 UG/KG	1.76
20	62	NOT FOUND							
21	117	1002	12:31	21	1.000	A BB	245709.	50.000 UG/KG	1.00
22	130	NOT FOUND							
23	63	659	8:14	9	1.084	A VB	29817.	17.206 UG/KG	0.34
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	804	10:03	21	0.802	A BB	154427.	86.918 UG/KG	1.74
27	92	812	10:09	21	0.810	A BB	2565720.	622.229 UG/KG	12.44
28	75	850	10:37	9	1.398	A BB	14331.	8.212 UG/KG	0.16
29	97	NOT FOUND							
30	164	NOT FOUND							
31	43	921	11:31	21	0.919	A VB	1185630.	838.206 UG/KG	16.75
32	129	NOT FOUND							
33	112	NOT FOUND							
34	106	1033	12:55	21	1.031	A BV	1557220.	575.321 UG/KG	11.50
35	106	1056	13:12	21	1.054	A BB	2145270.	531.968 UG/KG	10.64
36	106	1109	13:52	21	1.107	A BB	2051480.	568.526 UG/KG	11.37
37	104	1113	13:55	21	1.111	A BB	1068810.	164.268 UG/KG	3.28
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	539	6:44	1	1.164	A BB	109159.	48.513 UG/KG	0.97
41	95	1185	14:49	21	1.183	A VB	194920.	46.721 UG/KG	0.93
42	98	802	10:01	21	0.800	A BB	251153.	44.065 UG/KG	0.88

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:55	0.98	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:07		10.000			50.00		0.930	
3	1:13		10.000			50.00		1.260	
4	1:31		10.000			50.00		1.632	
5	1:37		10.000			50.00		0.710	
6	2:40		5.000			50.00		1.612	
7	2:50		5.000			50.00		4.085	
8	2:52	0.97	10.000	0.05	249.87	50.00	1.565	0.313	5.00
9	7:42	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	3:28	0.98	5.000	0.12	599.34	50.00	21.806	1.819	11.99
11	3:54		5.000			50.00		1.643	
12	4:36		5.000			50.00		2.402	
13	4:52		10.000			50.00		0.538	
14	5:34		5.000			50.00		1.673	

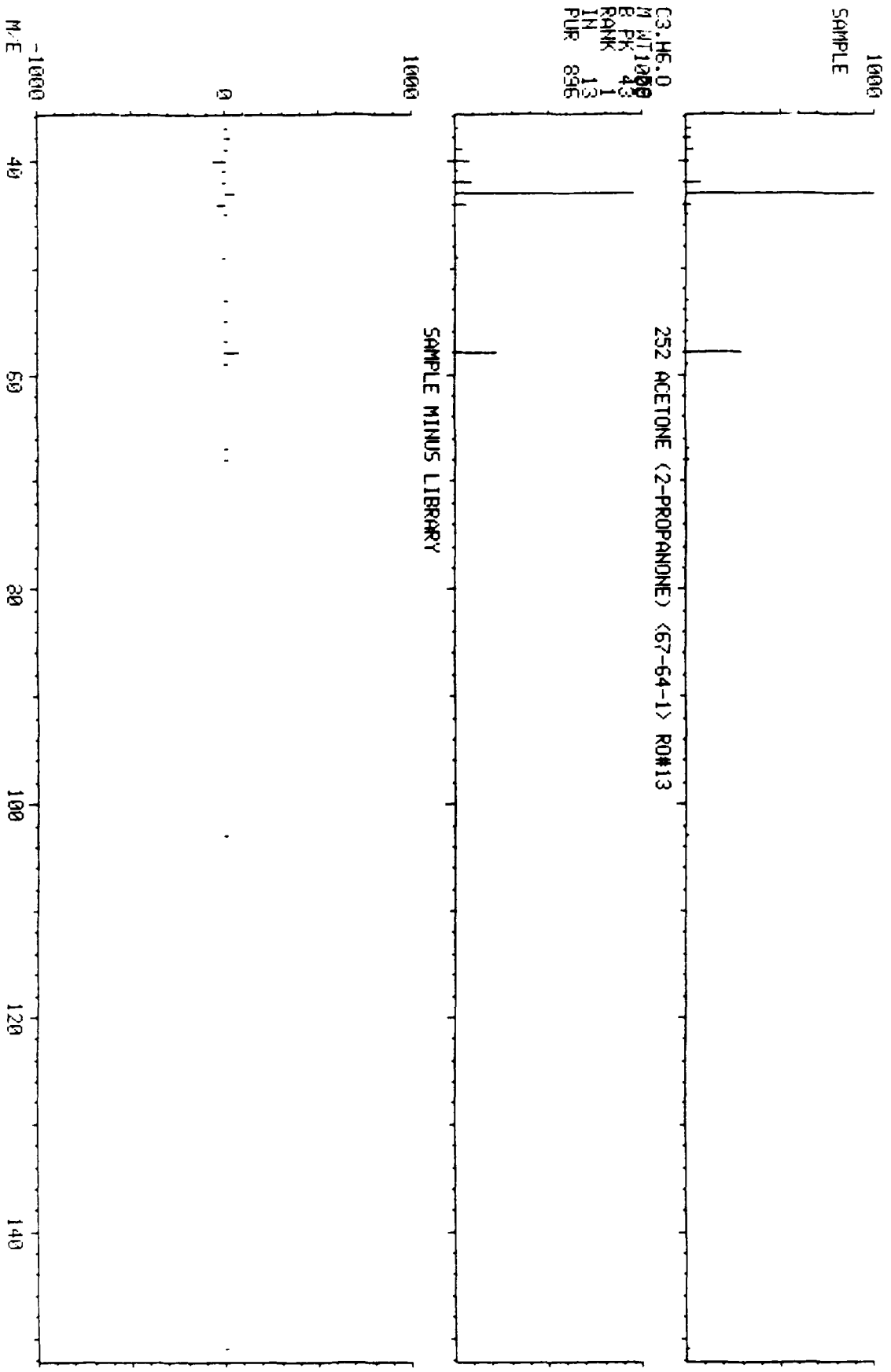
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:41	0.98	10.000	0.10	362.15	50.00	0.975	0.135	7.24
16	6:10		5.000			50.00		3.249	
17	6:19		5.000			50.00		0.880	
18	6:34		5.000			50.00		0.898	
19	6:53	0.99	5.000	0.18	88.10	50.00	1.756	0.997	1.76
20	6:58		5.000			50.00		2.119	
21	12:34	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	8:01		5.000			50.00		0.581	
23	8:21	0.99	5.000	0.22	17.21	50.00	0.122	0.354	0.34
24	8:55		5.000			50.00		0.785	
25	9:41		5.000			50.00		0.771	
26	10:04	1.00	15.000	0.05	86.92	50.00	0.628	0.362	1.74
27	10:10	1.00	5.000	0.16	622.23	50.00	10.442	0.839	12.44
28	10:41	0.99	5.000	0.28	8.21	50.00	0.059	0.357	0.16
29	10:57		5.000			50.00		0.407	
30	11:06		5.000			50.00		0.673	
31	11:34	1.00	15.000	0.06	838.21	50.00	4.825	0.288	16.76
32	11:34		5.000			50.00		0.776	
33	12:37		5.000			50.00		1.212	
34	12:55	1.00	5.000	0.21	575.32	50.00	6.338	0.551	11.51
35	13:09	1.00	5.000	0.21	531.97	50.00	8.731	0.821	10.64
36	13:51	1.00	5.000	0.22	568.53	50.00	8.349	0.734	11.37
37	13:54	1.00	5.000	0.22	164.27	50.00	4.350	1.324	3.29
38	14:07		5.000			50.00		0.577	
39	15:15		5.000			50.00		0.655	
40	6:51	0.98	5.000	0.23	48.51	50.00	1.604	1.653	0.97
41	14:49	1.00	5.000	0.24	46.72	50.00	0.793	0.849	0.93
42	10:04	1.00	5.000	0.16	44.07	50.00	1.022	1.160	0.88

COMPUchem LABS
LIBRARY SEARCH
12/27/89 7:18:00 + 2:47
SAMPLE: 1G CC#809687 CASE#18756.7 EPA#B202TAR ON#19
ENHANCED (5 158 24 01)

DATA: GR009687019 # 223

BASE M/E: 43
RIC: 19967.

C3.H6.O
M WT 100.0
E PK 43
RANK 1
IN 13
PUR 896

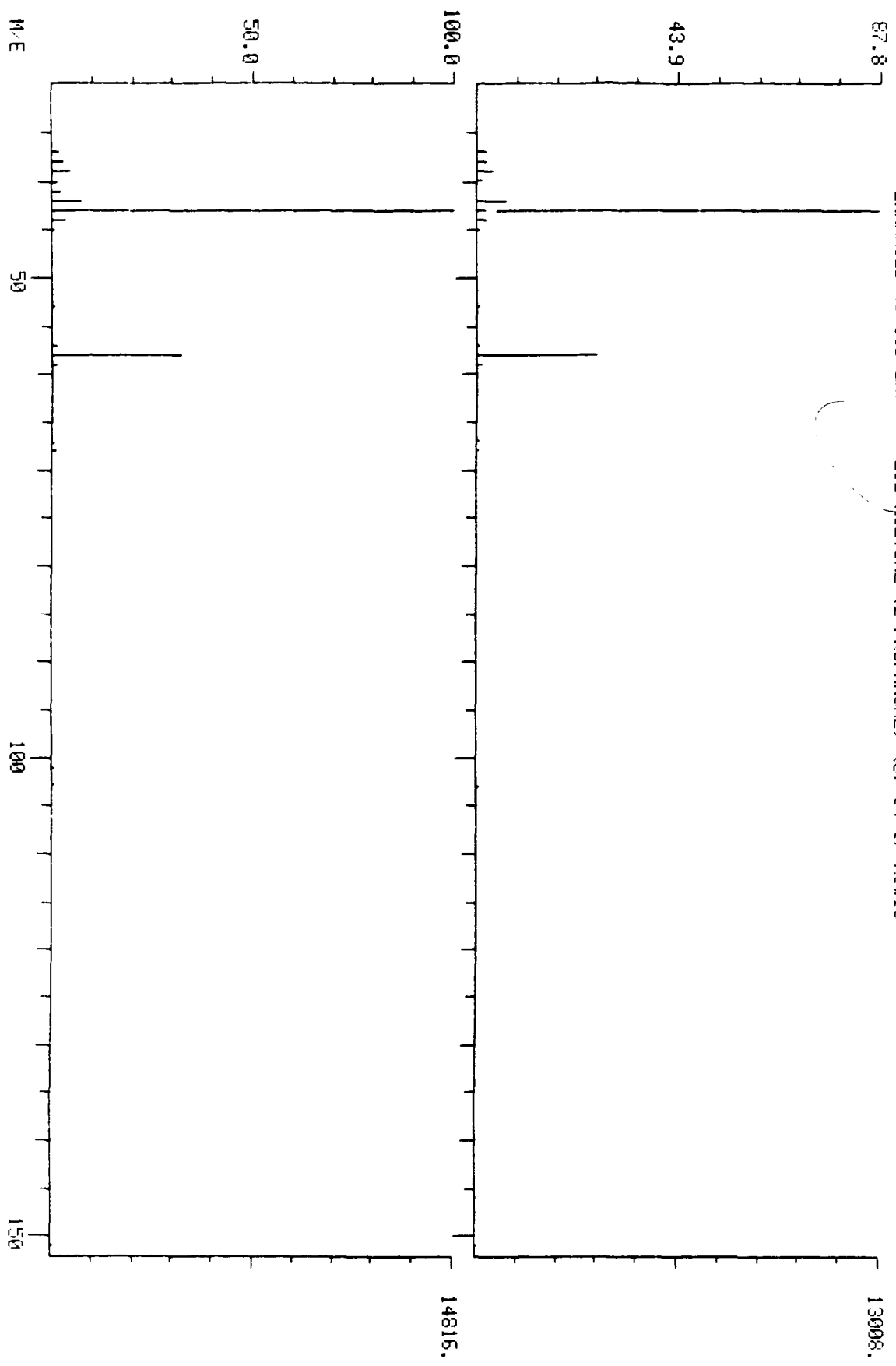


QUAL MASS SPECTRUM
12/27/69 7:10:00 + 2:47
SAMPLE: 1G.00#309697 CASE#18756.7 EPA#B202TAR.0N#19
ENHANCED (S 158.2N) 252 ACETONE (2-PROPANONE) (67-64-1) R0#13

COMPUHER LABS

DATA: GR009697C19 #223

BASE M/E: 43 53
RIC: 19967. 23807.



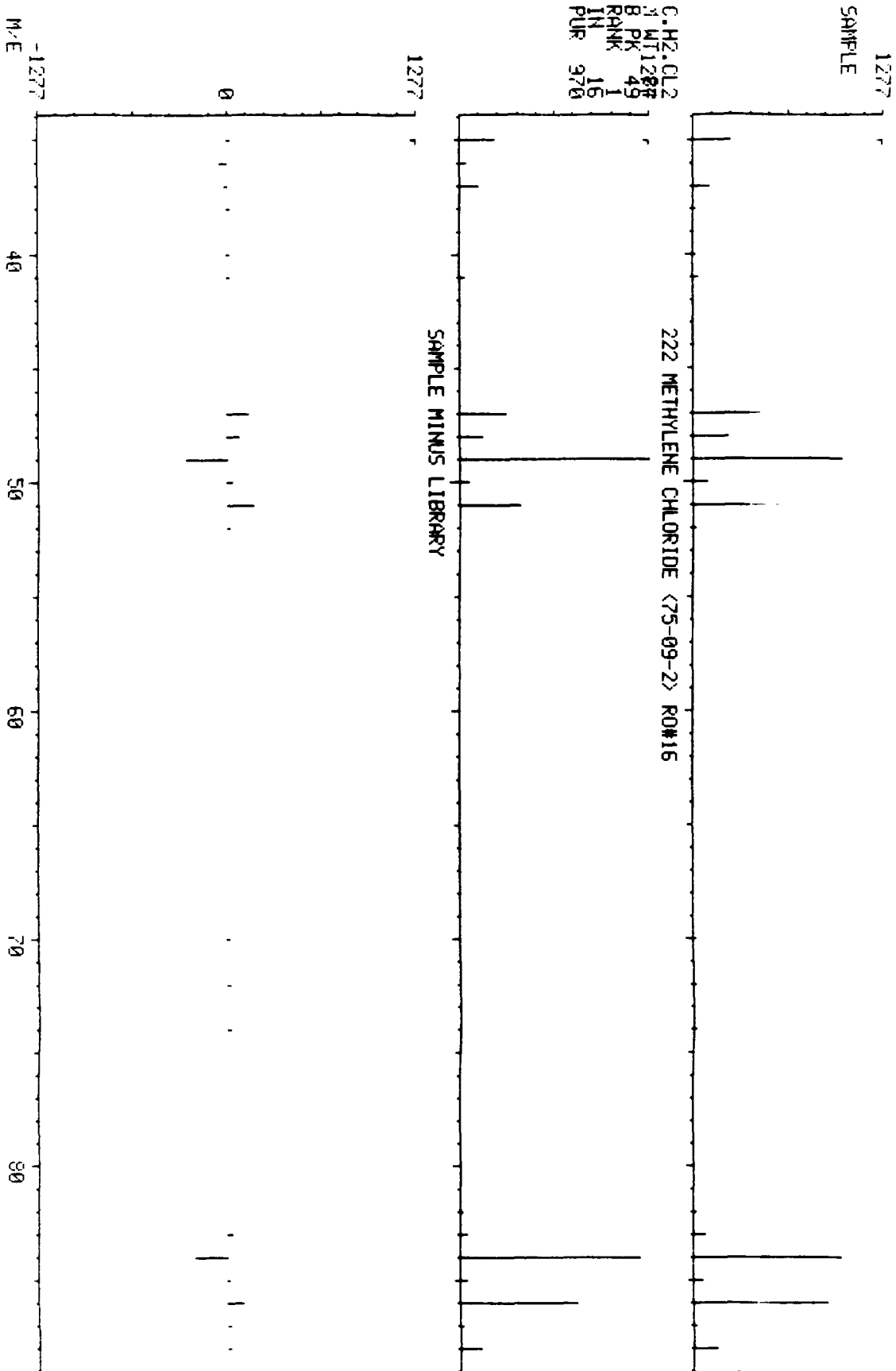
LIBRARY SEARCH
12/27/89 7:18:00 + 3:24
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B2021AR QN#19
ENHANCED (5 158 2H 0T)

COMPUCHEM LABS

DATA: GR009637C19 # 272

BASE M/E: 84
RIC: 655359.

C.H2.CL2
X WT 12.87
B PK 49
RANK 1
IN 16
PUR 970



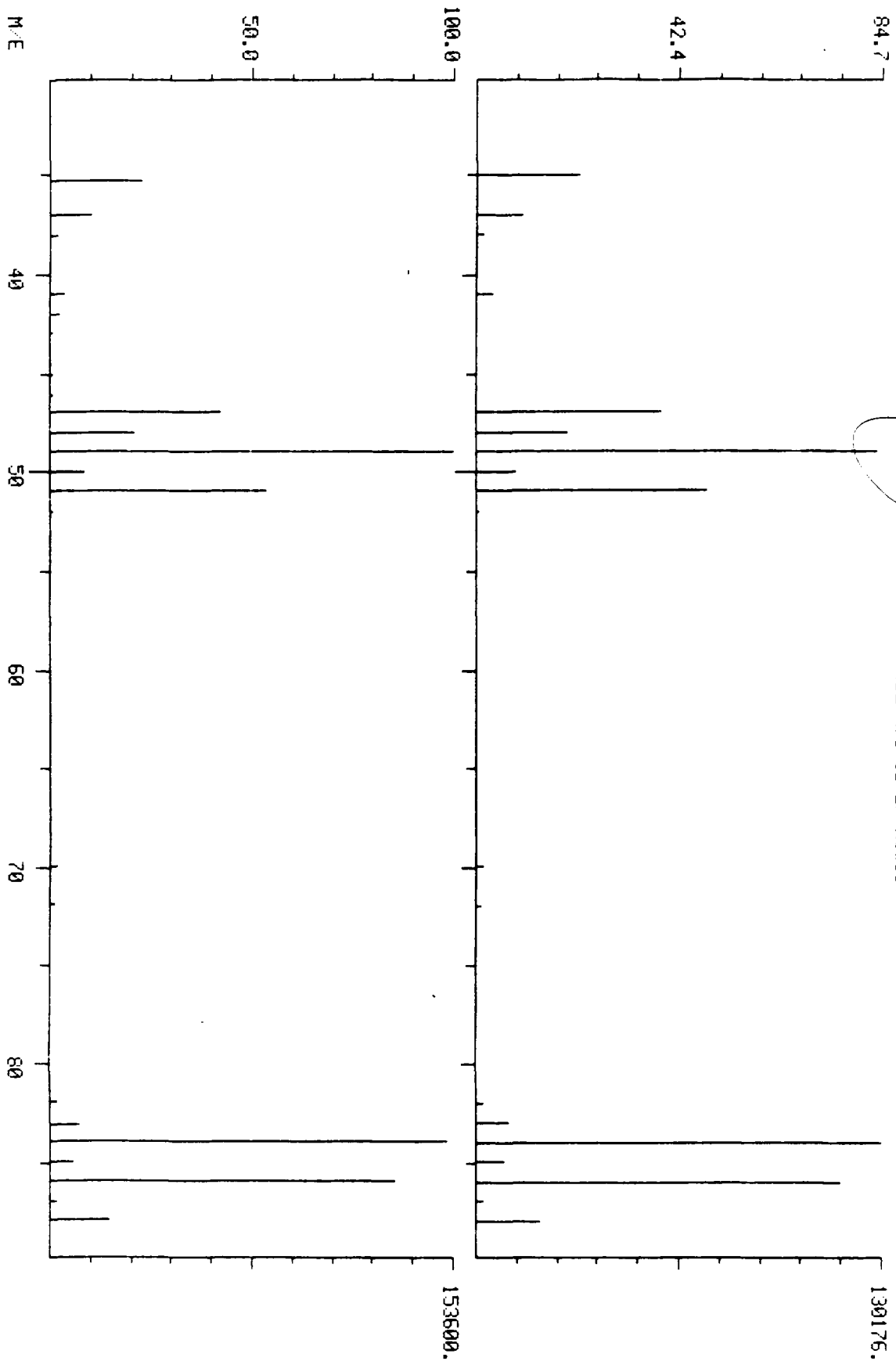
DUAL MASS SPECTRUM
12/27/89 7:10:00 + 3:24
SAMPLE: 1G OC#309687
ENHANCED (S 158 2M)

18756.7 EPA#B202TAR ON#19
222 METHYLENE CHLORIDE <75-09-2> R0#16

COMPUCHEN LABS

DATA: GR009687C19 #272

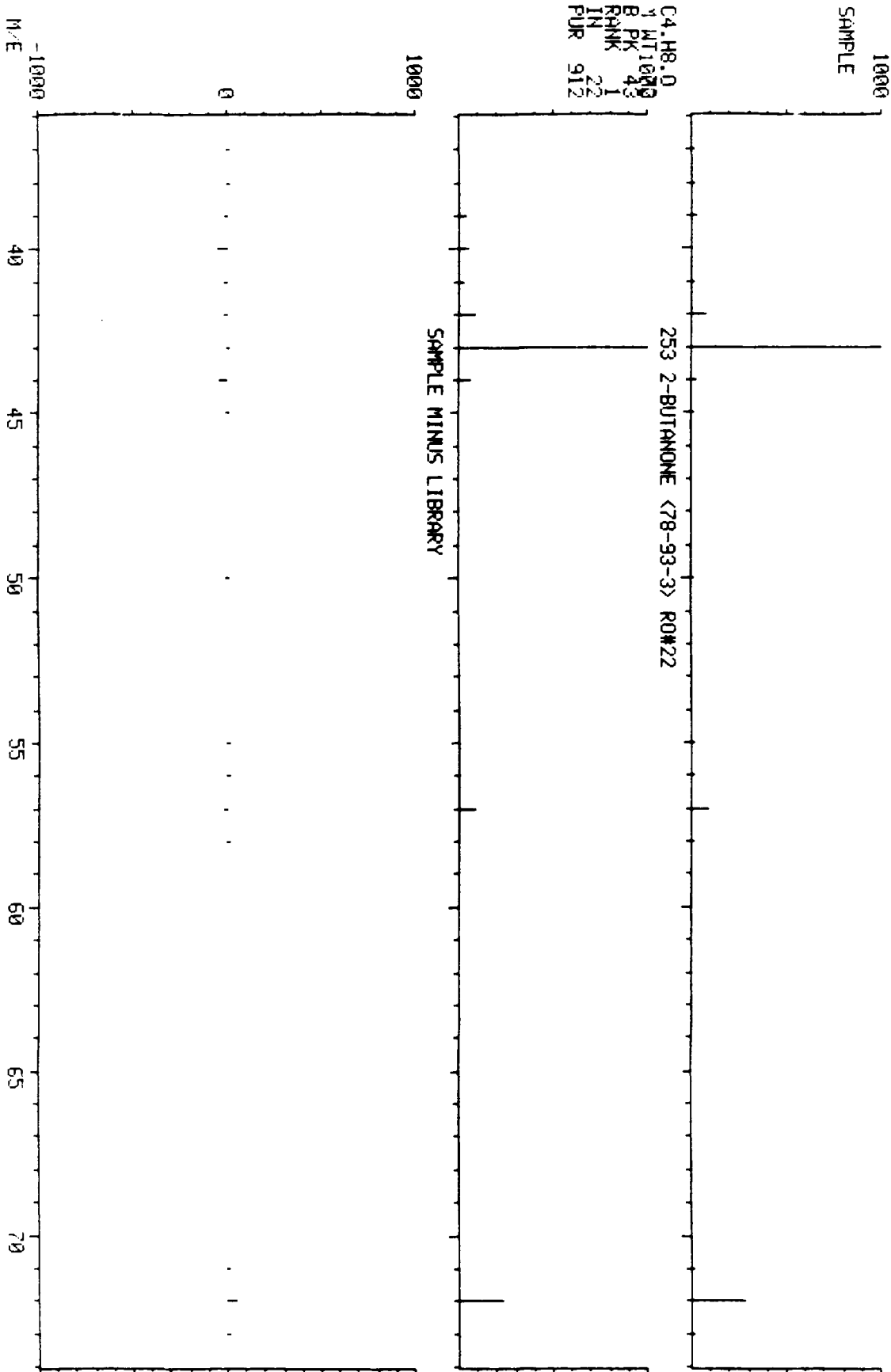
BASE M/E: 84/ 49
RIC: 655359. / 746495.



COMPUCHEN LABS
LIBRARY SEARCH
12/27/89 7:10:00 + 5:34
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR QN#19
ENHANCED (5.156 ZN QT)

DATA: GR09687019 # 445
BASE M/E: 43
RIC: 37567.

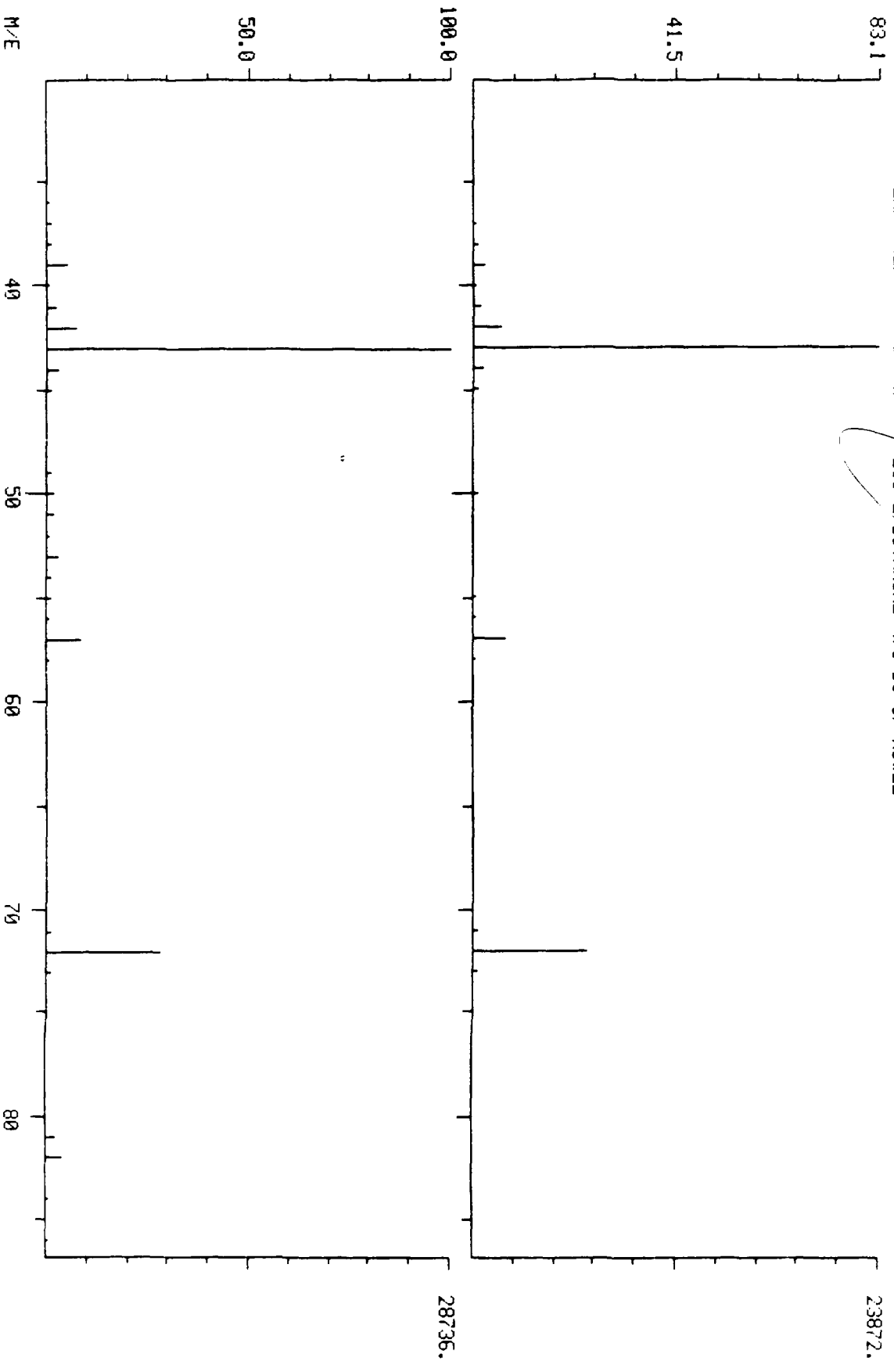
C4.H8.O
M WT 100.0
B PK 43
RANK 1
IN 22
PUR 912



DUAL MASS SPECTRUM
12/27/89 7:10:00 + 5:34
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B2021AR QN#19
ENHANCED (5 15B 2N) 253 2/BUTANONE (78-93-3) RM#22

CONFLUENCE LABS

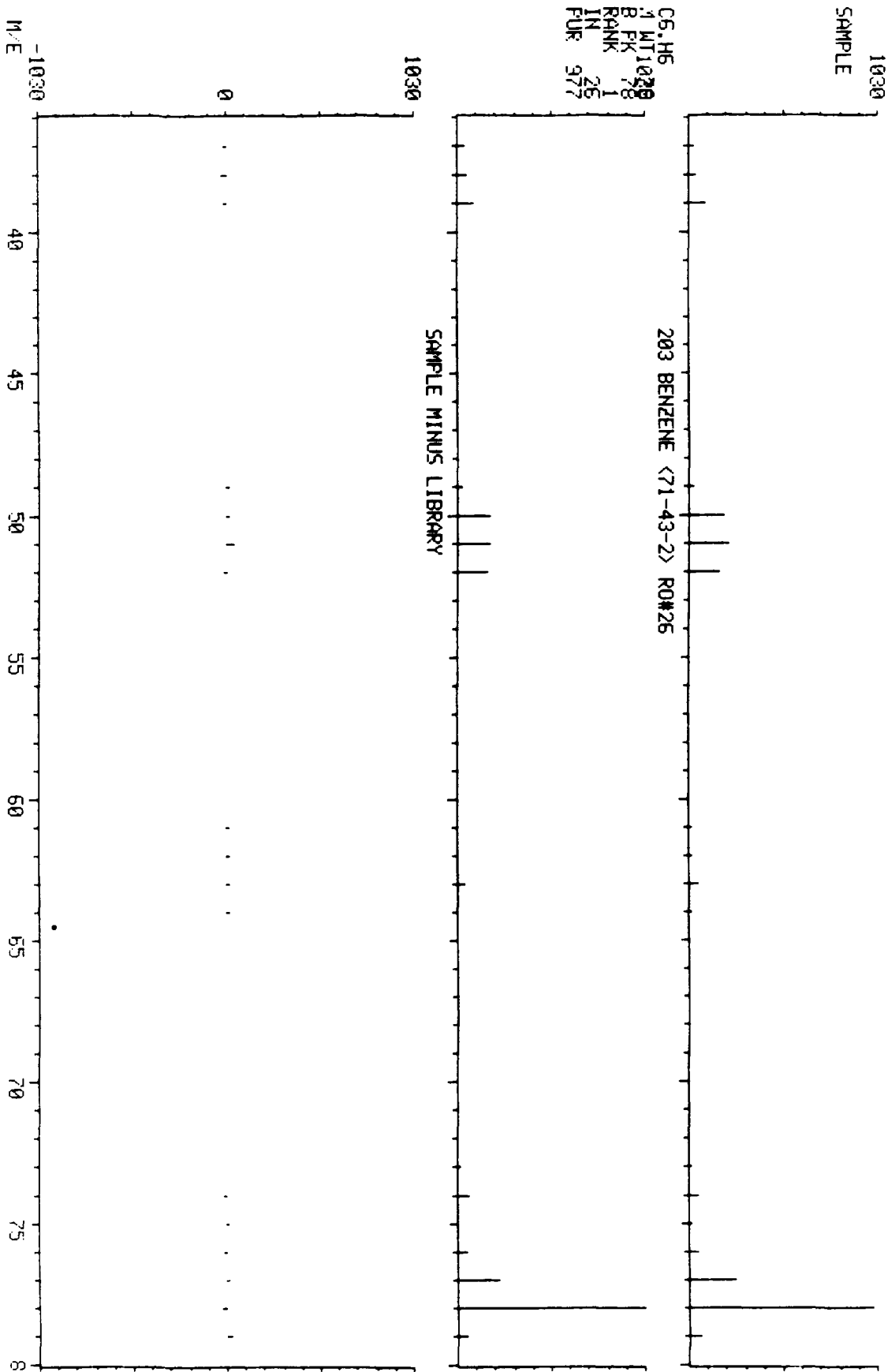
DATA: GR009687C19 #445 BASE M/E: 43/ 43
RIC: 37695./ 51391.



LIBRARY SEARCH
12/27/89 7:10:00 + 5:47
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR QM#19
ENHANCED (5 158 2H 0T)

COMPILED: LAES
DATA: GR009687019 # 543
BASE M/E: 78
R/C: 87935.

06.H6
1 WT 1020
8 PK 78
RANK 1
IN 26
PUR 977

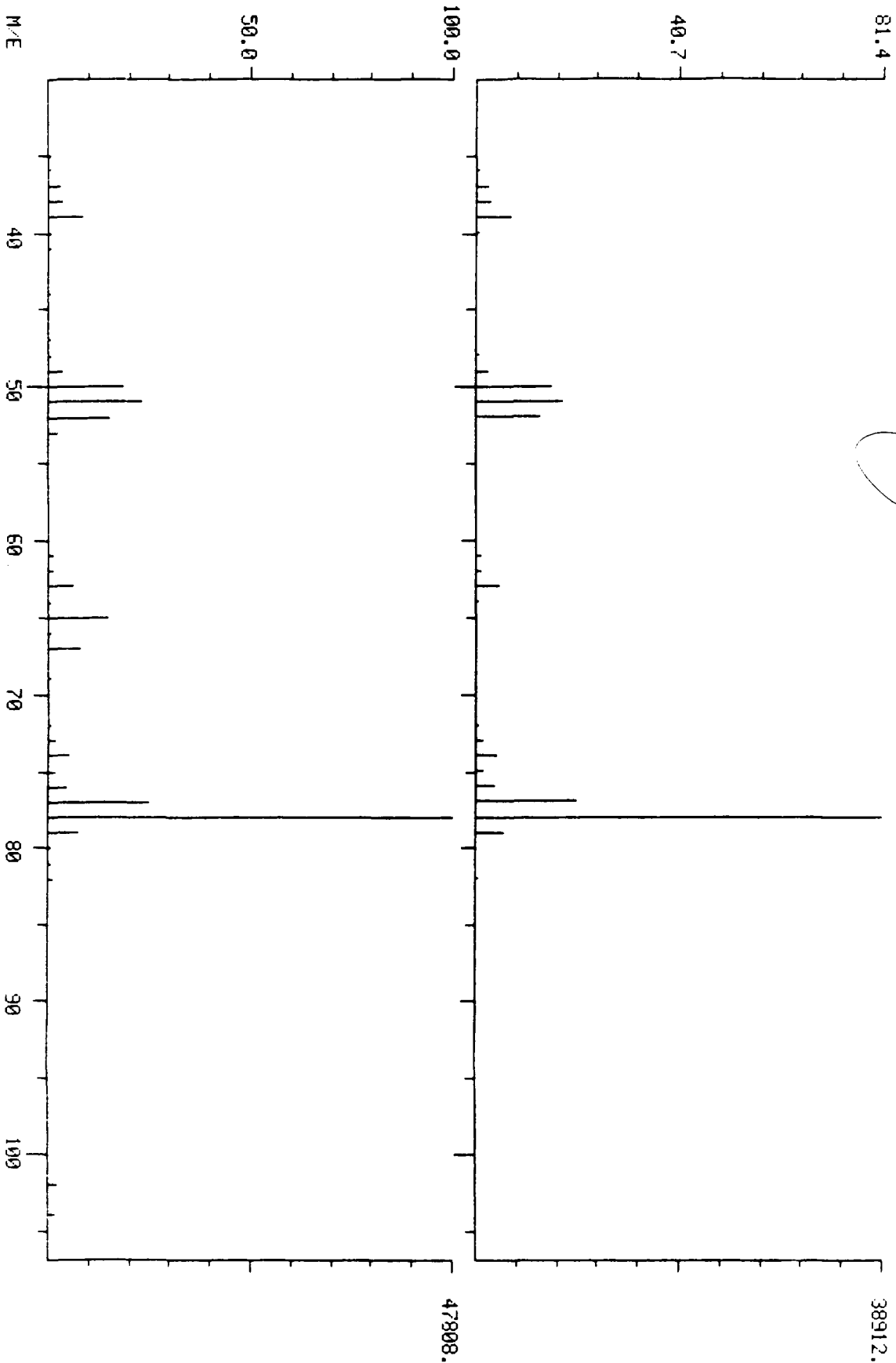


DUAL MASS SPECTRUM
12/27/89 7:10:00 + 6:47
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR ON#19
ENHANCED (5 15B 2N) 203 BENZENE (71-43-2) RM#26

COMFUCHEM LABS

DATA: GR009687C19 #543 BASE M/E: 78/ 78

RIC: 88959.7 126847.

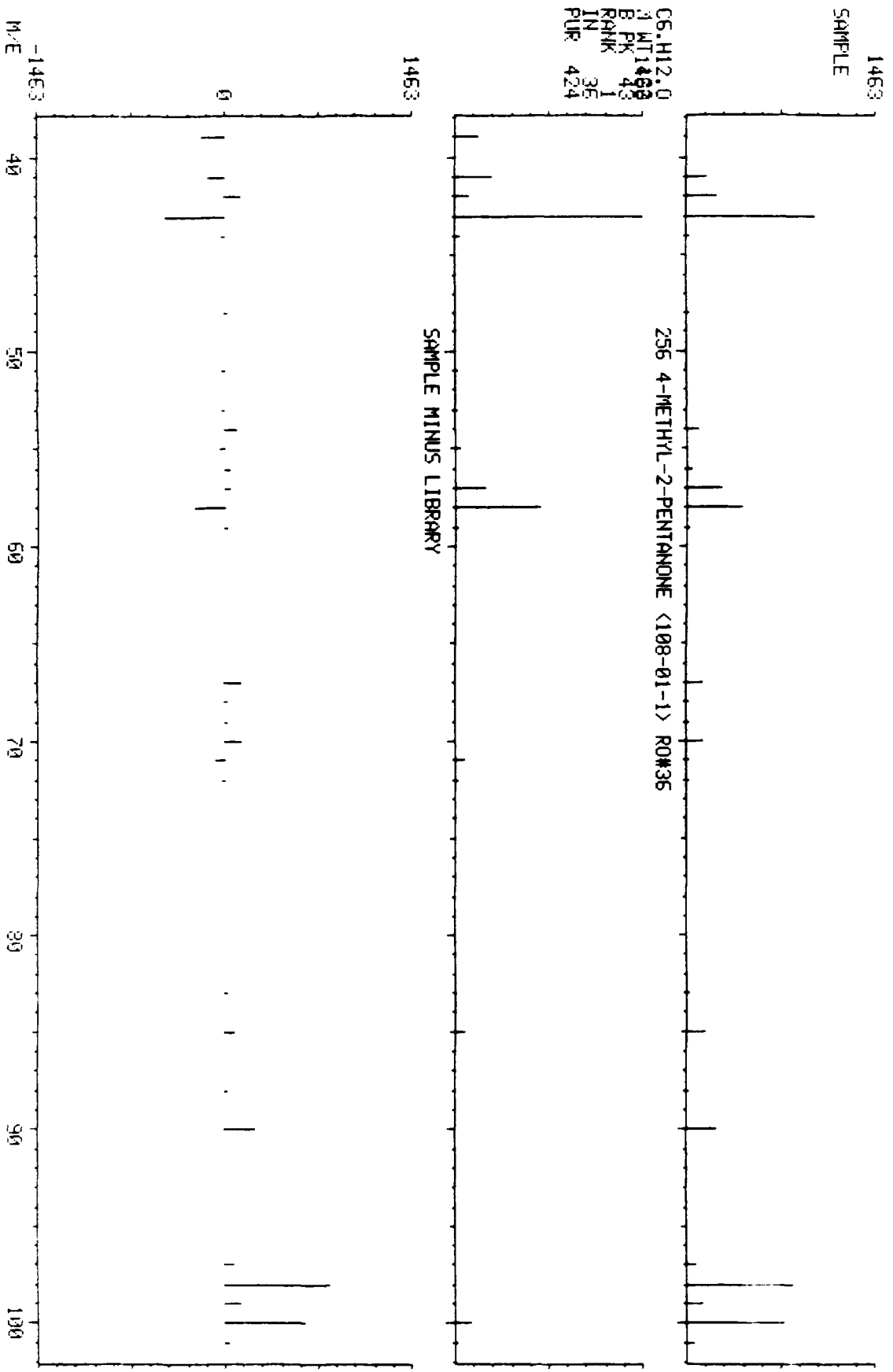


CONFUCHEM LABS
LIBRARY SEARCH
12/27/89 7:10:00 + 10:03
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR QN#19
ENHANCED (S 158 2H 0T)

DATA: GP005687C19 # 804

BASE M/E: 43
R10: 28719.

06.H12.0
1.0
1463
1
36
424



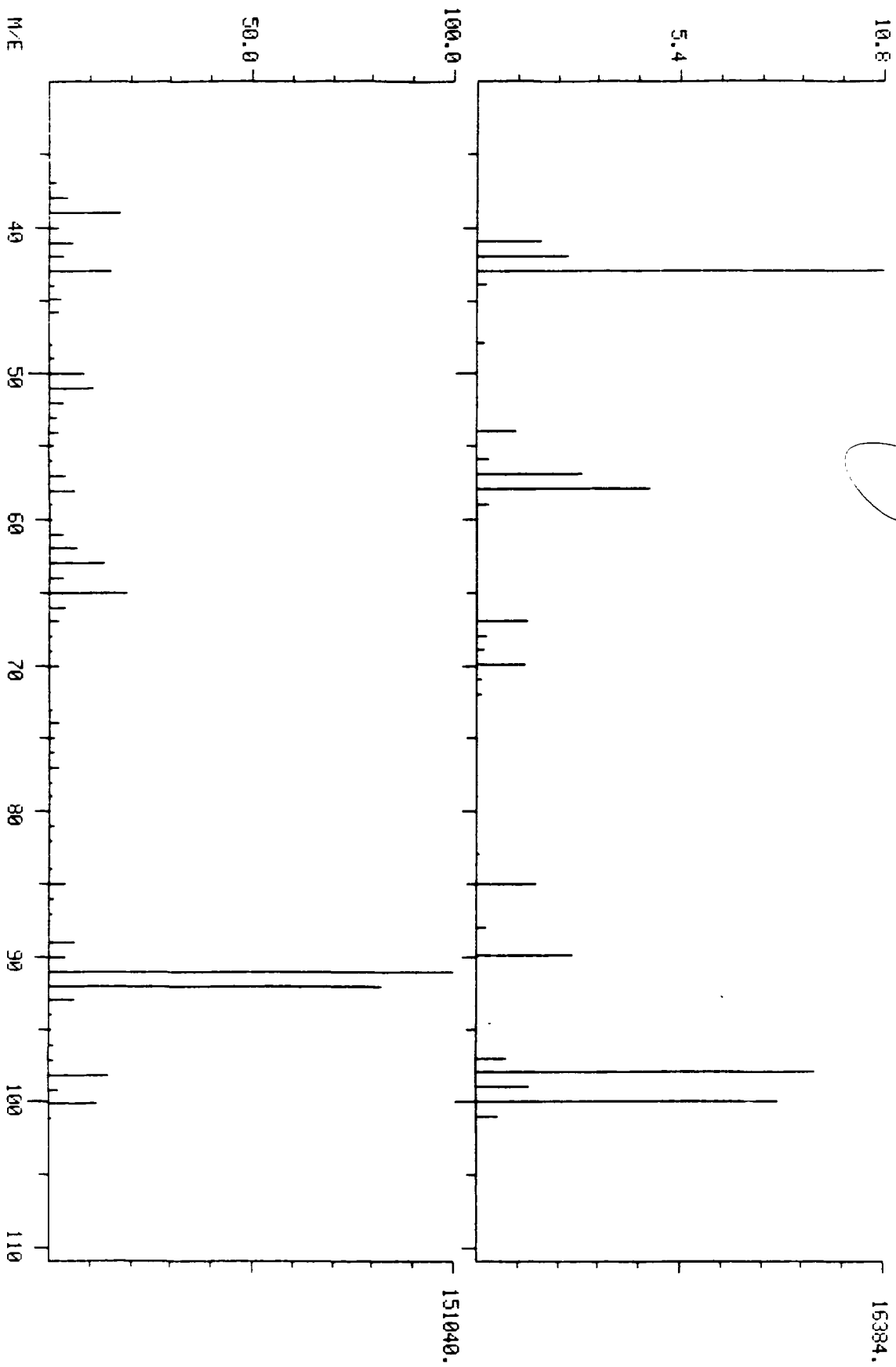
COMPUCHEM LABS

DATA: CR069687C19 #804

BASE M/E: 43/ 51

RIC: 78719.7 603135.

DUAL MASS SPECTRUM
12/27/89 7:10:00 + 10:03
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B2021AR QN#19
ENHANCED (5 158 2N) 256 4 METHYL-2-PENTANONE <108-01-1> R0#36

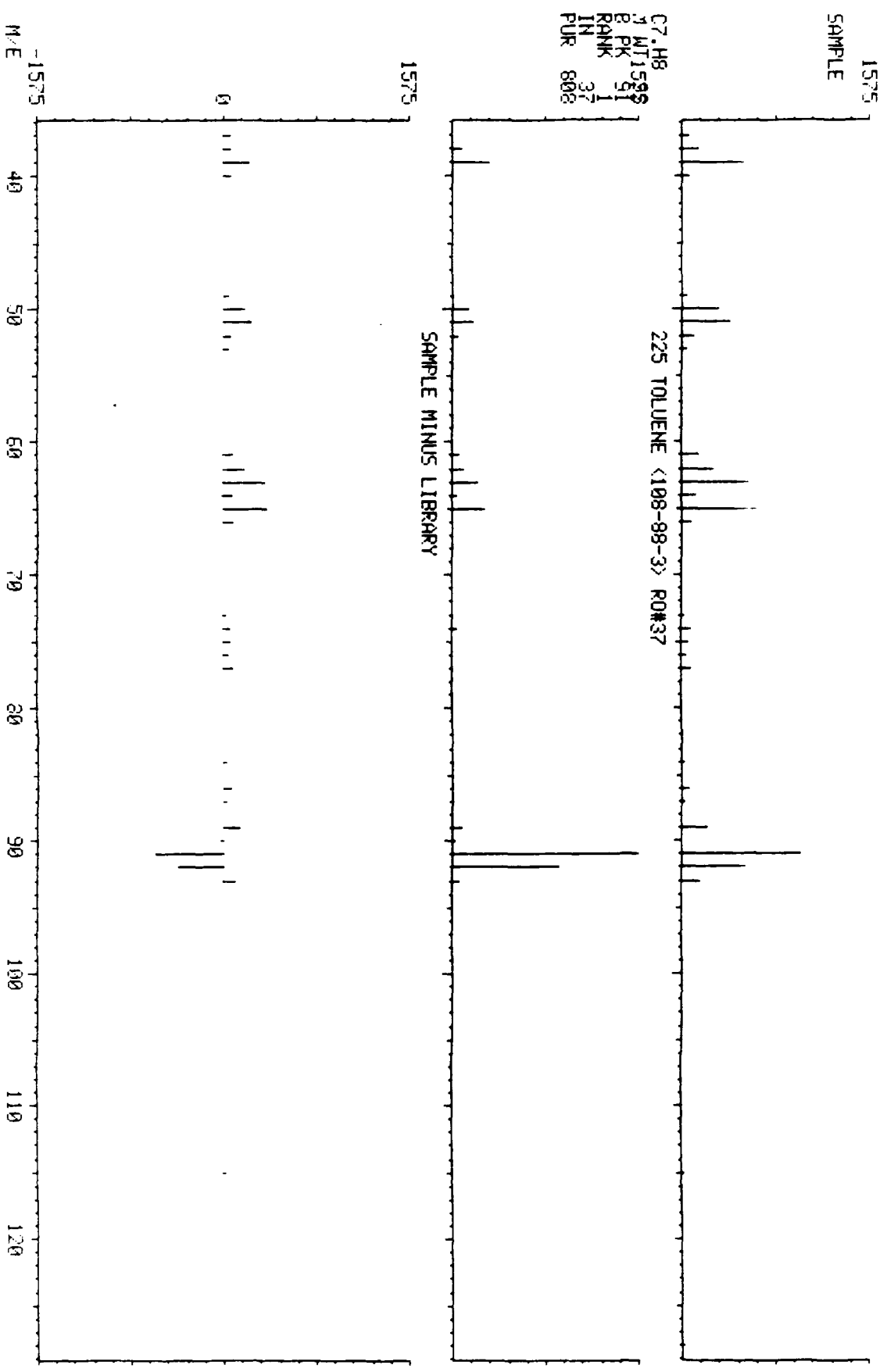


COMPUCHER LABS
LIBRARY SEARCH
12/27/89 7:10:00 + 10:09
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR ON#19
ENHANCED (5 158 2N 0T)

DATA: CR009687019 # 812

BASE M/E: 91
RIC: 549887.

C7.H8
7 WT 1595
8 PK 51
RANK 37
IN 37
PUR 800

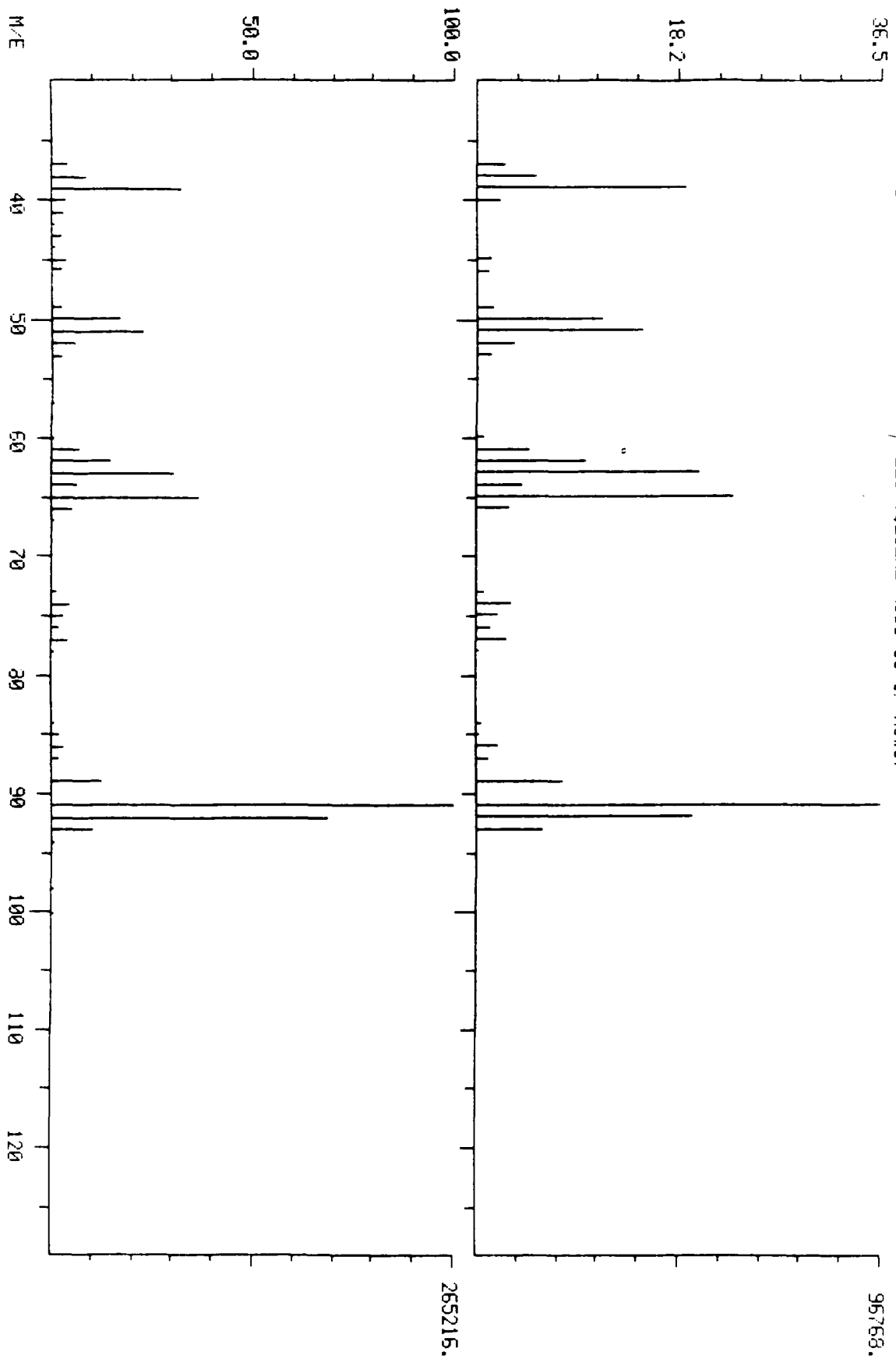


DUAL MASS SPECTRUM
12/27/89 7:10:00 + 18:03
SAMPLE: 1G CCM309687 CASE#18756.7 EPA#B202TAR QN#19
ENHANCED (5 158 2H) / 225 TOLUENE <108-88-3> R0#37

COMPUchem LABS

DATA: GR009687C19 #812

BASE M/E: 91 / 91
RIC: 559103. / 1128440.



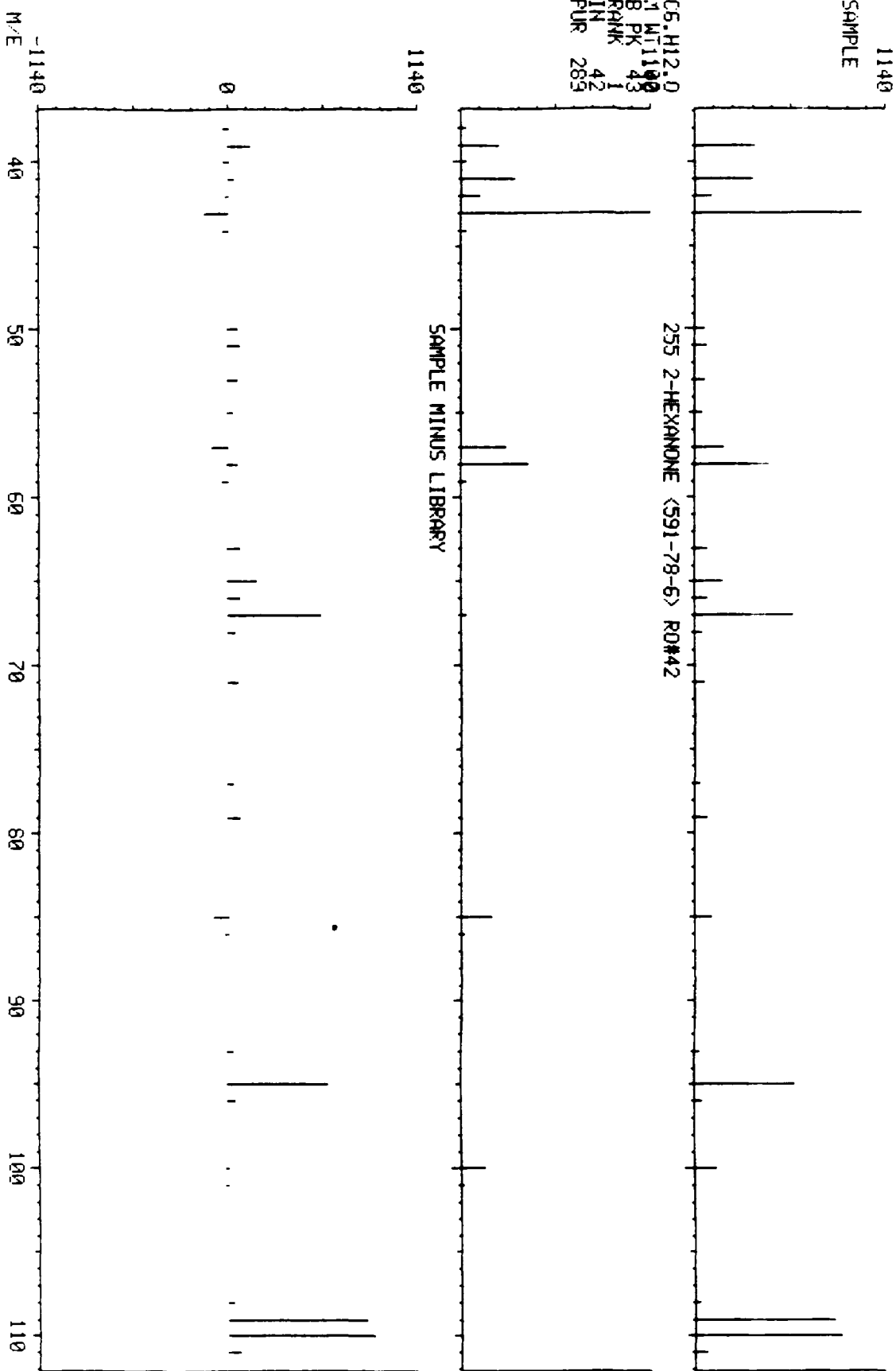
LIBRARY SEARCH
12/27/89 7:10:00 + 11:31
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B2021AR QN#19
ENHANCED (S 156 2N 0T)

COMPUCHEN LABS

DATA: GR009687C19 # 921

BASE M/E: 43
RIC: 550911.

06.H12.0
7.MI 119.0
8.PK 43
RANK 1
IN 42
PUR 283

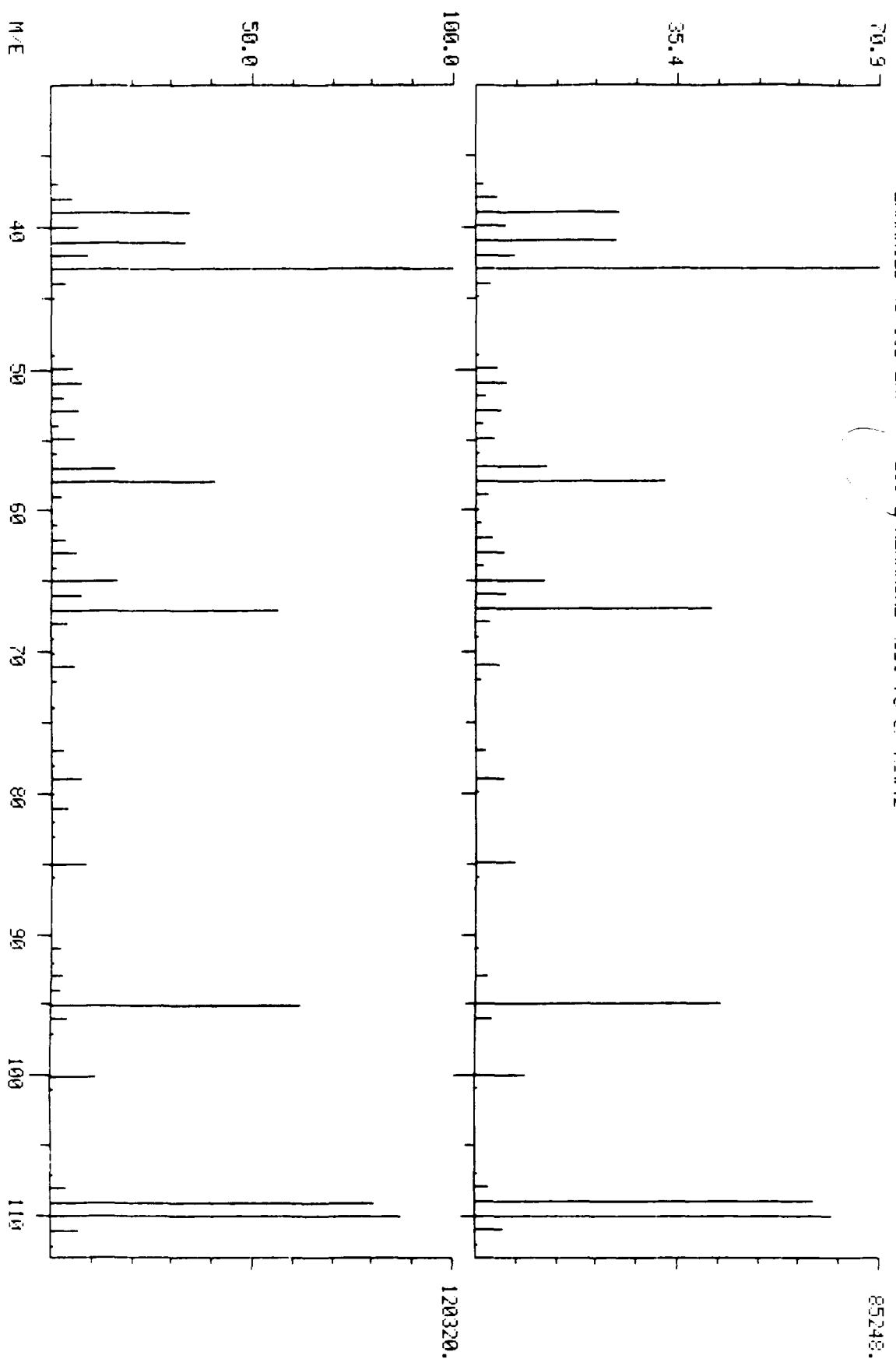


QUAL MASS SPECTRUM
12/27/89 7:10:00 + 11:31
SAMPLE: 16 CC#309687 CASE#18756.7 EPA#B202TAR OH#19
ENHANCED (S-158 2N) 255 2-HEXANONE (591-79-6) RM#42

COMPUCHEN LABS

DATA: 06069687019 #921

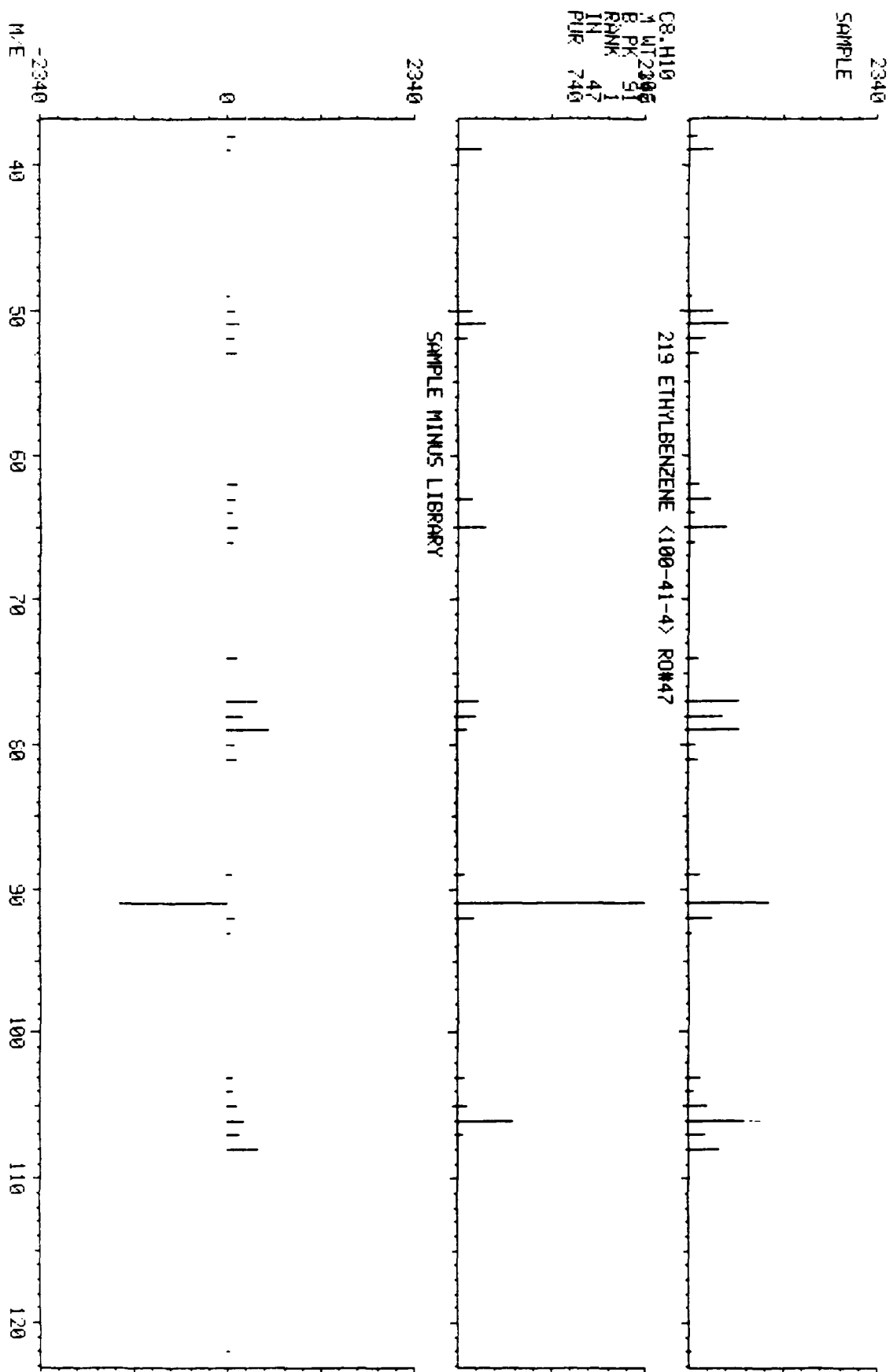
BASE M/E: 43/ 43
RID: 563579. / 218175.



COMPUCHEN LABS
 DATA: GR009687019 #1033
 BASE M/E: 91
 R1C: 685055

LIBRARY SEARCH
 12/27/89 7:10:00 + 12:55
 SAMPLE: 16 OC#309687 CASE#18756.7 EPA#B202TAR ON#19
 ENHANCED (5 158 2H 0T)

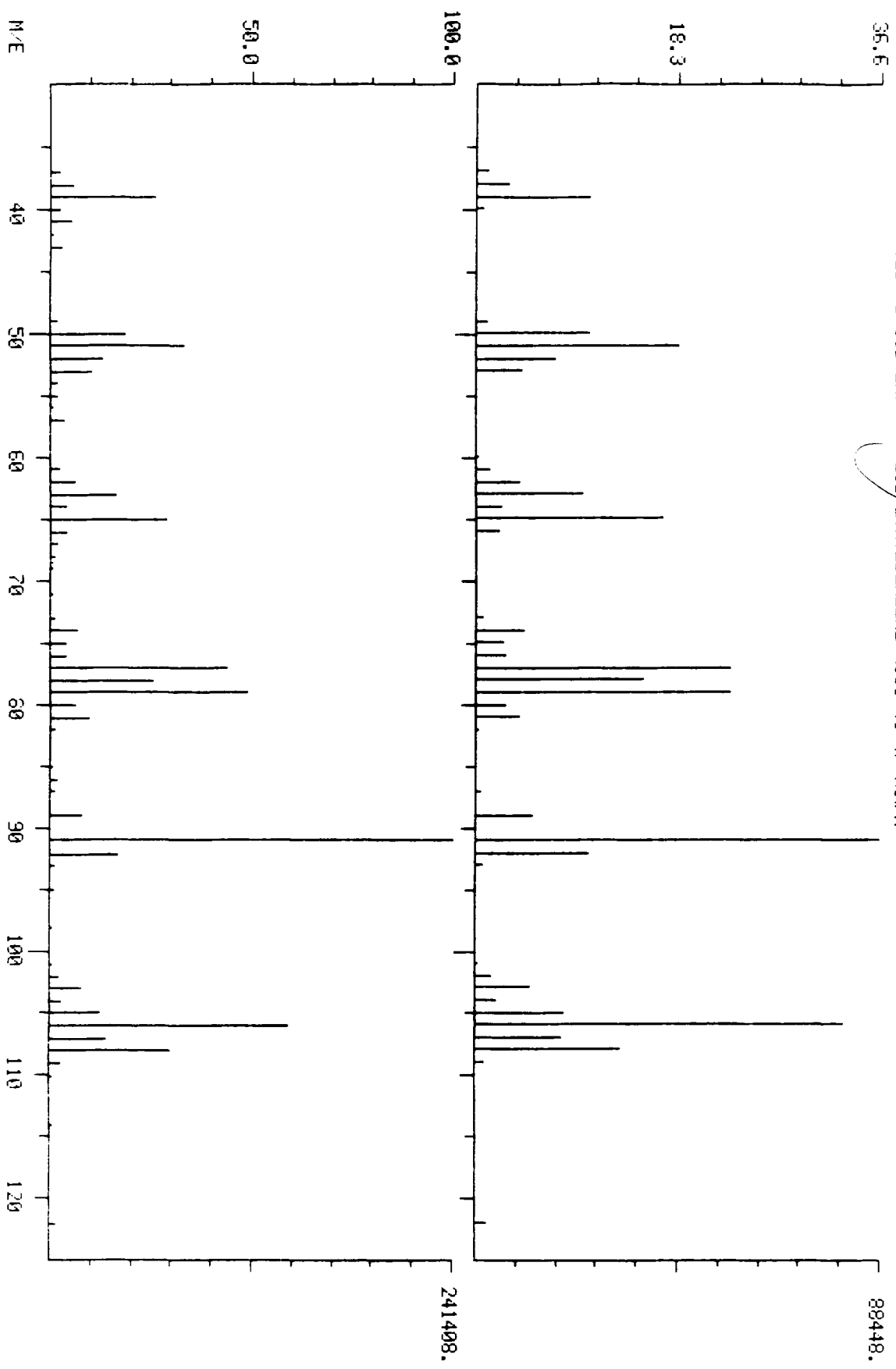
08.H10
 1 WT 206
 8 PK 51
 RANK 1
 IN 47
 PUR 740



DUAL MASS SPECTRUM
12/27/89 7:10:00 + 12:55
SAMPLE: 15 CC#309687 CASE#18756.7 EPA#R202TAR CH#19
ENHANCED (5 158.2N) 219/ETHYLENEZENE <100-41-4> F0#47

COMPUCHEN LABS

DATA: GR009687019 #1033 BASE M/E: 91, 91
RIC: 713727. / 1470460.



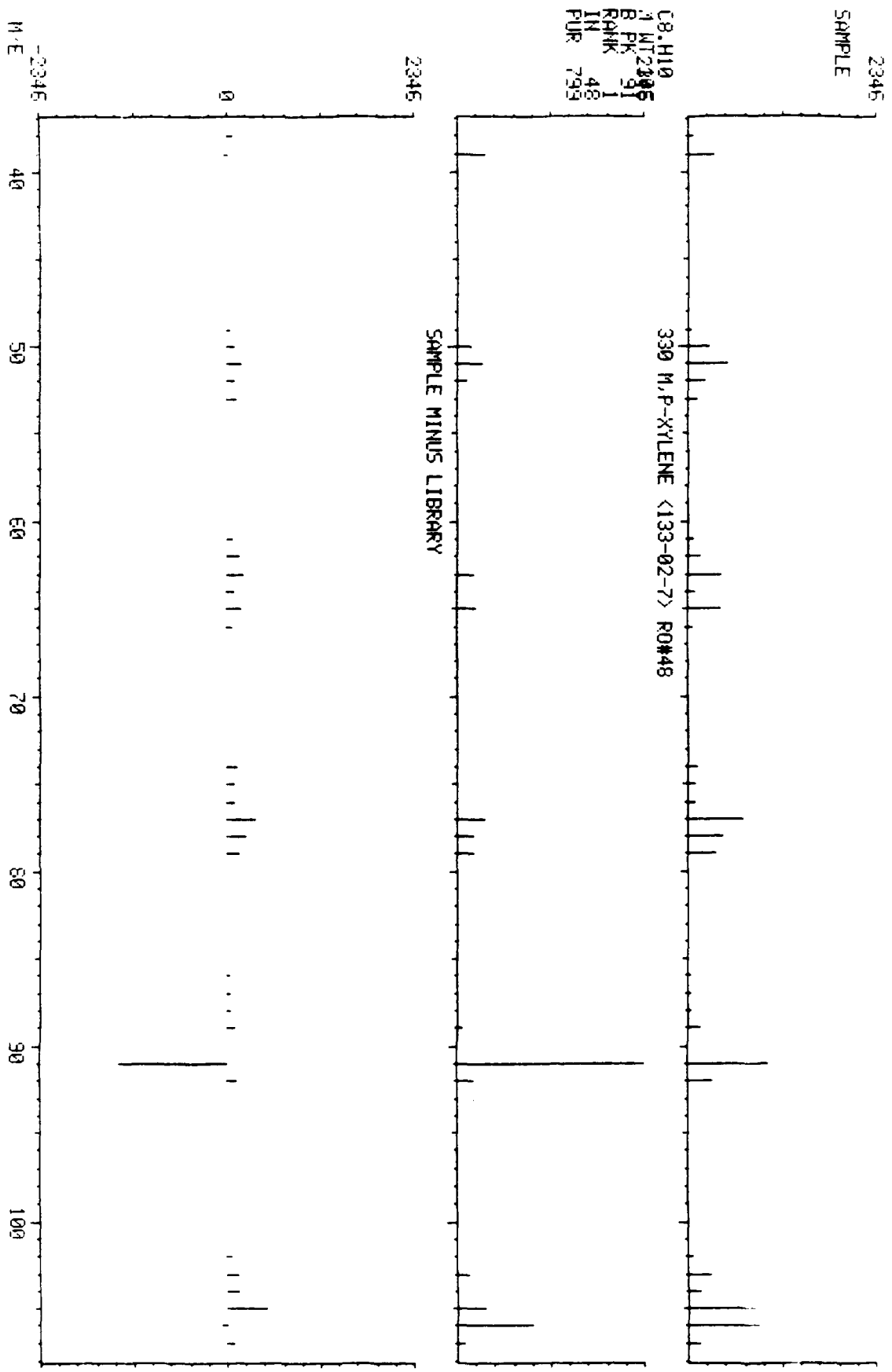
LIBRARY SEARCH
12/27/89 7:10:00 + 13:12
SAMPLE: 1G OC#399687 CASE#18756.7 EPA#B202TAR ON#19
ENHANCED (S 156 2N 0T)

COMPUCHEM LABS

DATA: GR009667C19 #1056

BASE M/E: 91
RIC: 802815.

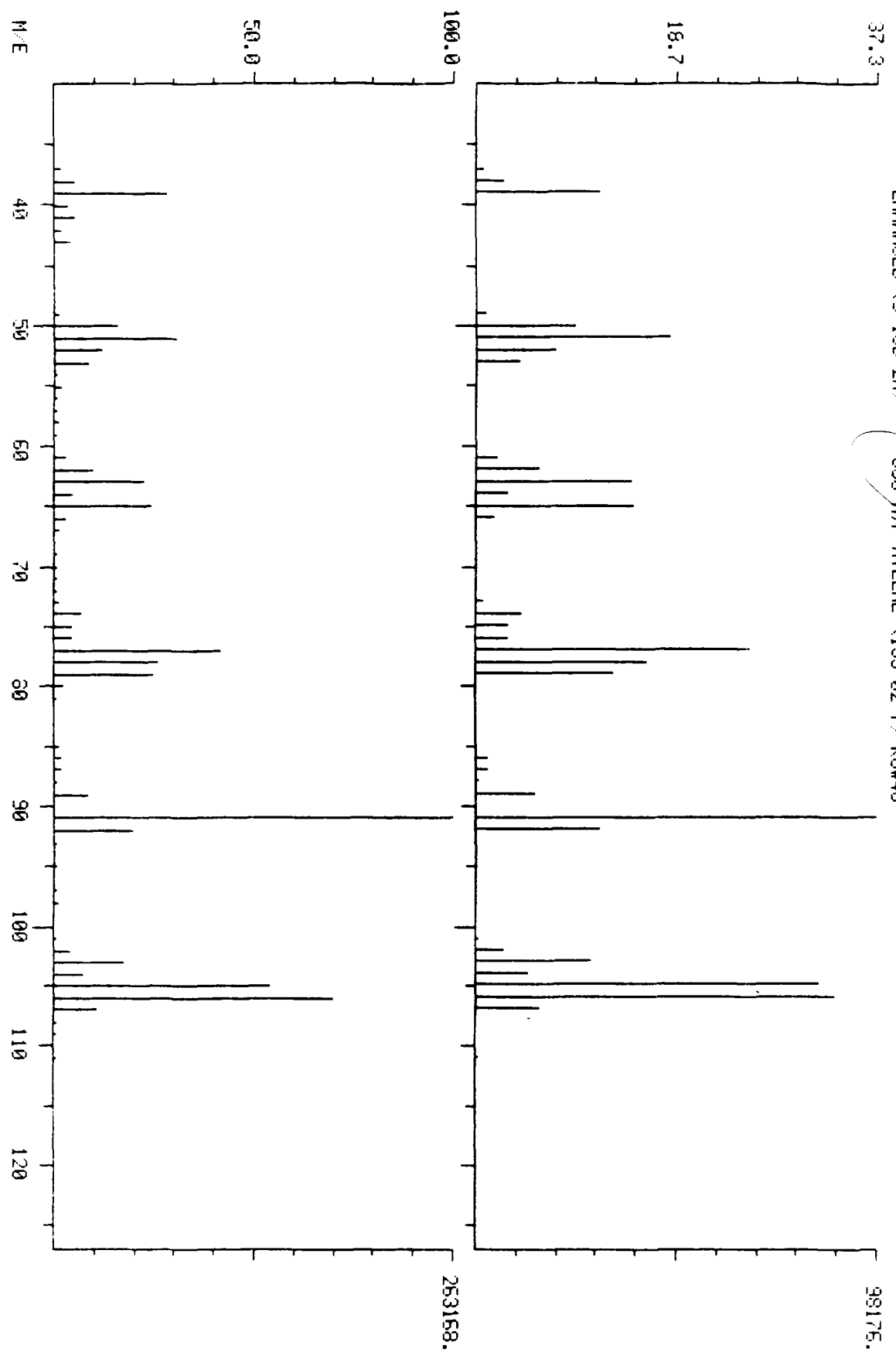
C8-H10
7 WITZ
B PK 91
RANK 1
IN 48
PUR 799



DUAL MASS SPECTRUM
12/27/89 7:10:00 + 13:12
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR ON#19
ENHANCED (5 158 ZN) 330 M.P.-XYLENE (133-02-7) RD#48

COMPUCHEN LABS

DATA: GR009687019 #1056 BASE M/E: 91/ 91
RIC: 808959.7 1597430.



COMPUCHER LABS

DATA: GR009687019 #1109

BASE M/E: 91

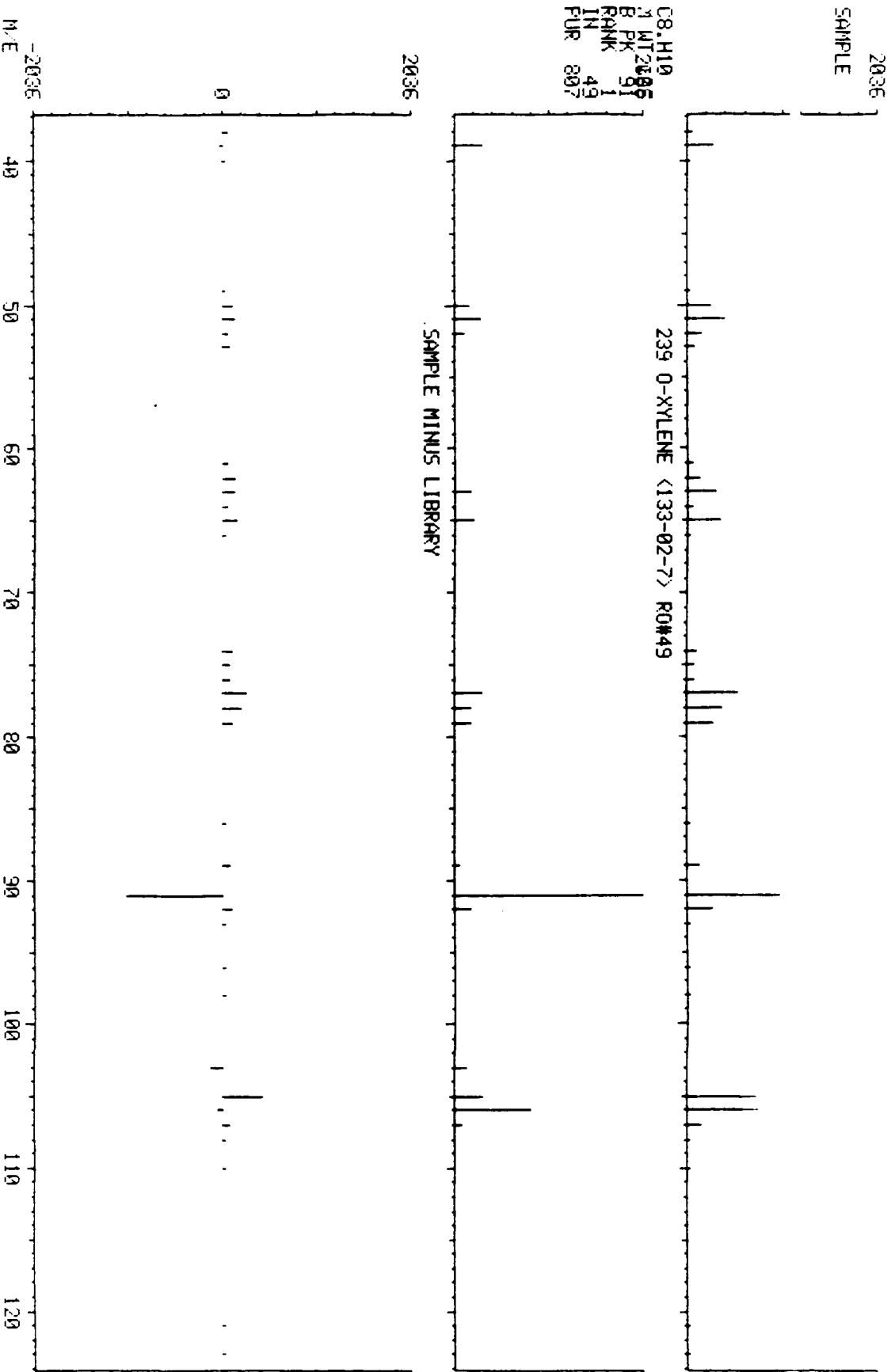
RIC: 615423.

LIBRARY SEARCH

12/27/89 7:10:00 + 13:52

SAMPLE: 1G (C#309687 CASE#18756.7 EPA#B202TAR QN#19

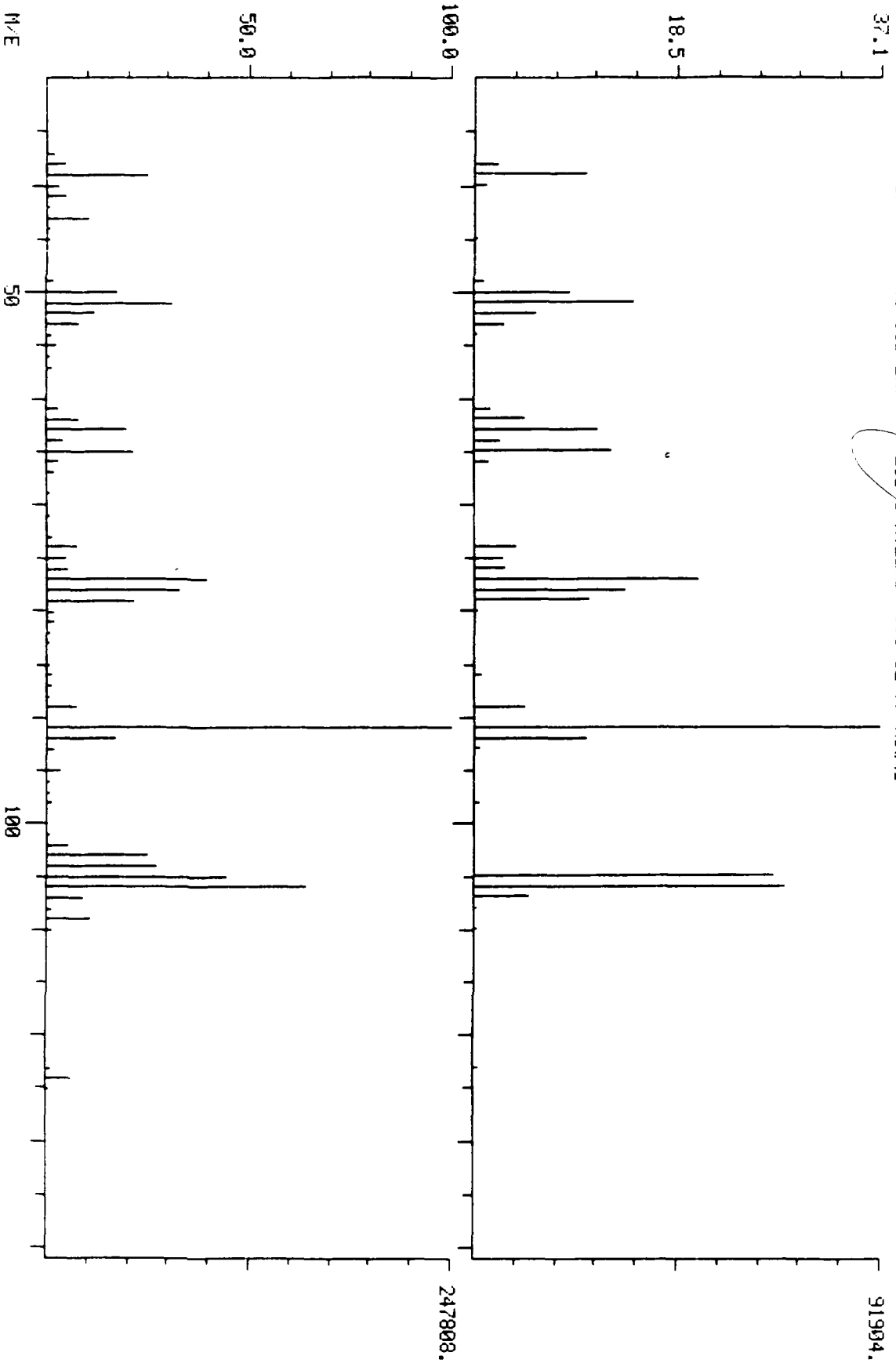
ENHANCED (5.156 2H 0T)



DUAL MASS SPECTRUM
12/27/89 7:10:00 + 13:52
SAMPLE: 16 CC#309687 CASE#18756.7 EPAMB202TAR QH#19
ENHANCED (5 158 2N) 239 0-XYLENE (133-02-7) RM#49

COMPUCHEN LABS

DATA: GR069687C15 #1109 BASE M/E: 91 / 91
RIC: 616447. / 1576950.

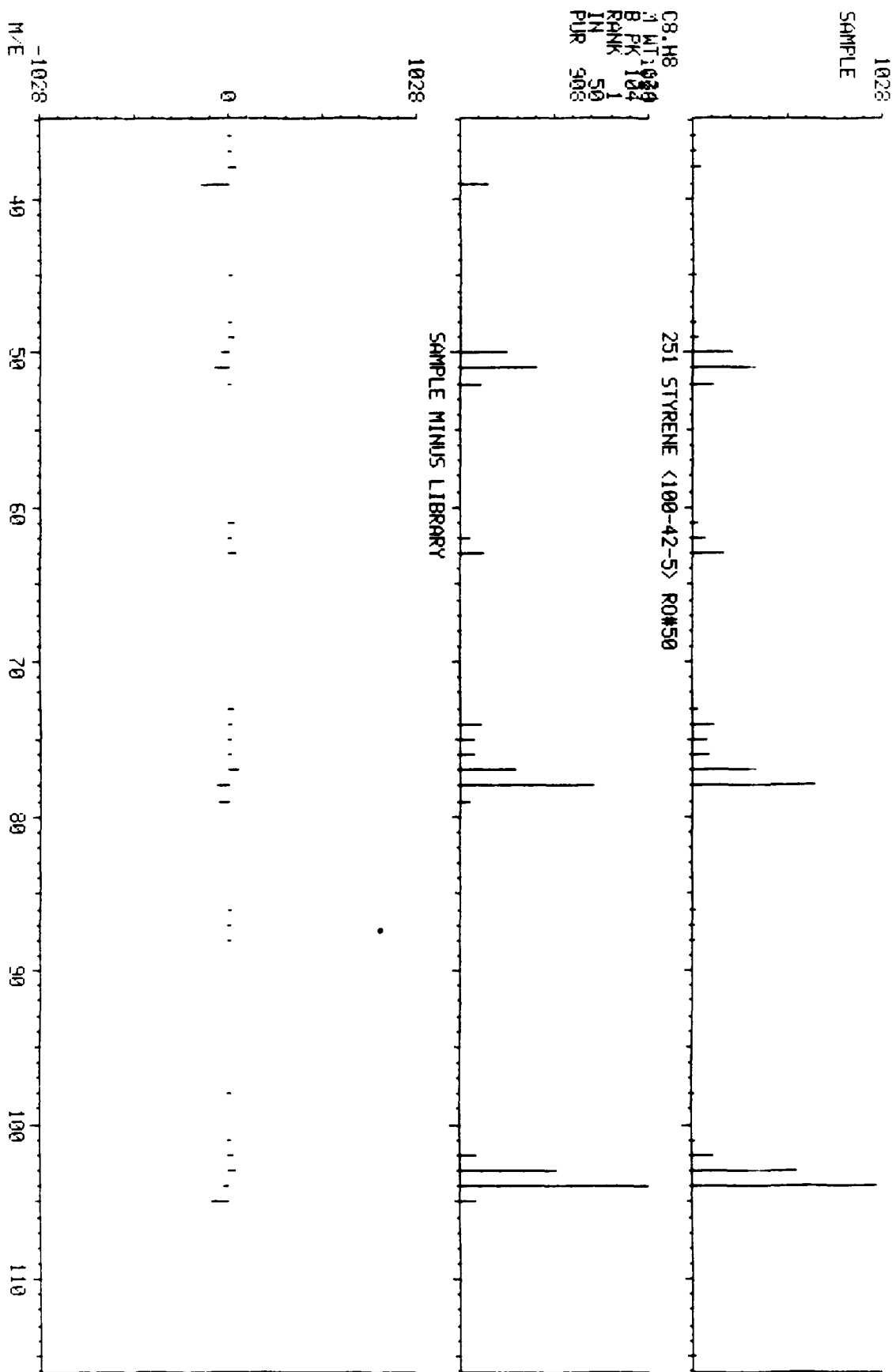


LIBRARY SEARCH
12/27/89 7:10:00 + 13:55
SAMPLE: 1G CC#309687 CASE#18756.7 EPAMB202TAP QM#19
ENHANCED (5 158 2N QT)

COMPUTER LABS

DATA: GR009687019 #1113

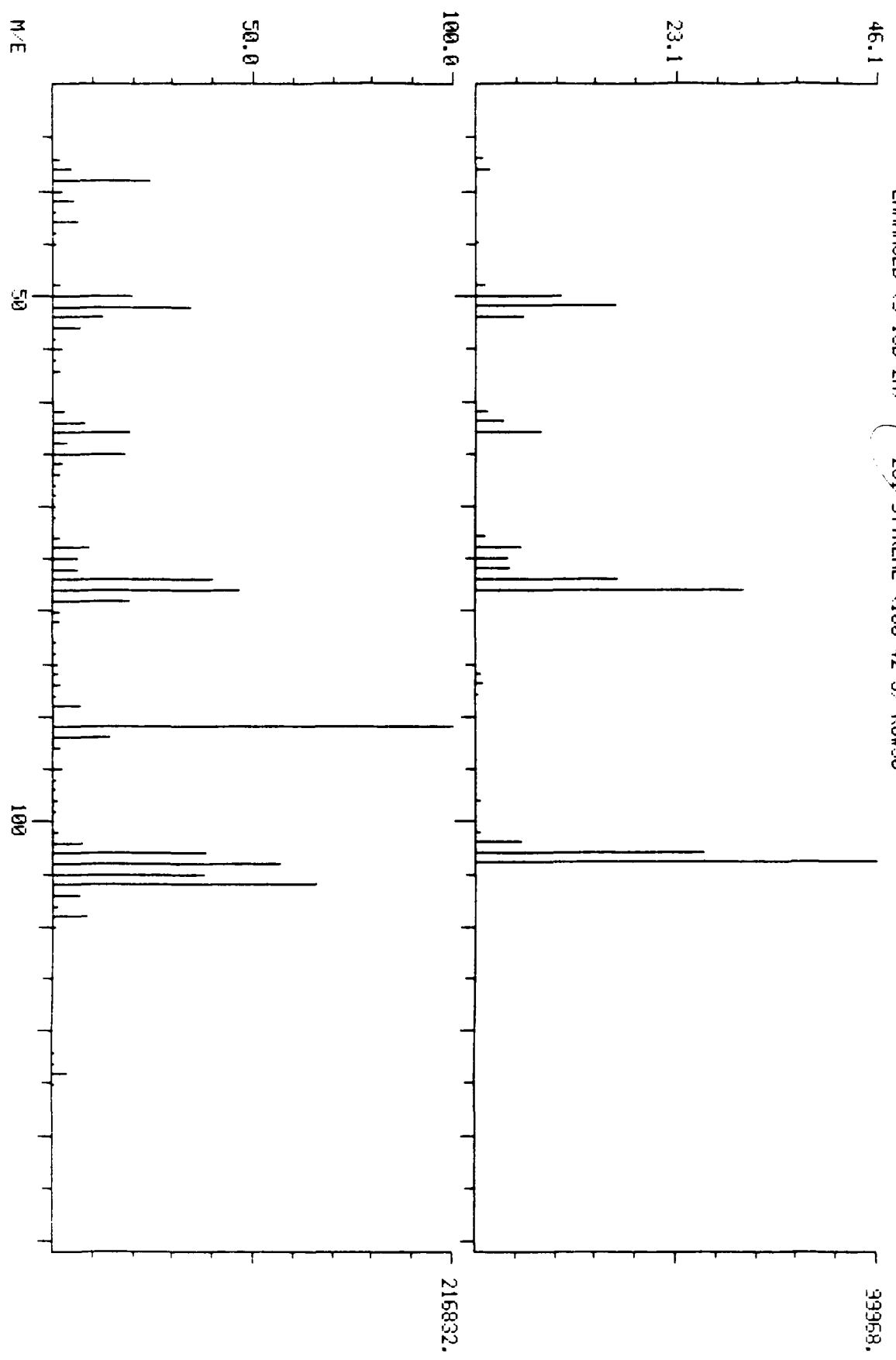
BASE M/E: 104
RIC: 408575.



DUAL MASS SPECTRUM
12/27/89 7:10:00 + 13:55
SAMPLE: 1G.CC#309687 CASE#18756.7 EPA#E202TAR.QM#19
ENHANCED (S 158.2M) 251 STYRENE <100-42-5> RM#50

COMPUCHEM LABS

DATA: GR009687C19 #1113 BASE M/E: 104 / 91
RIC: 400063. / 1462410.

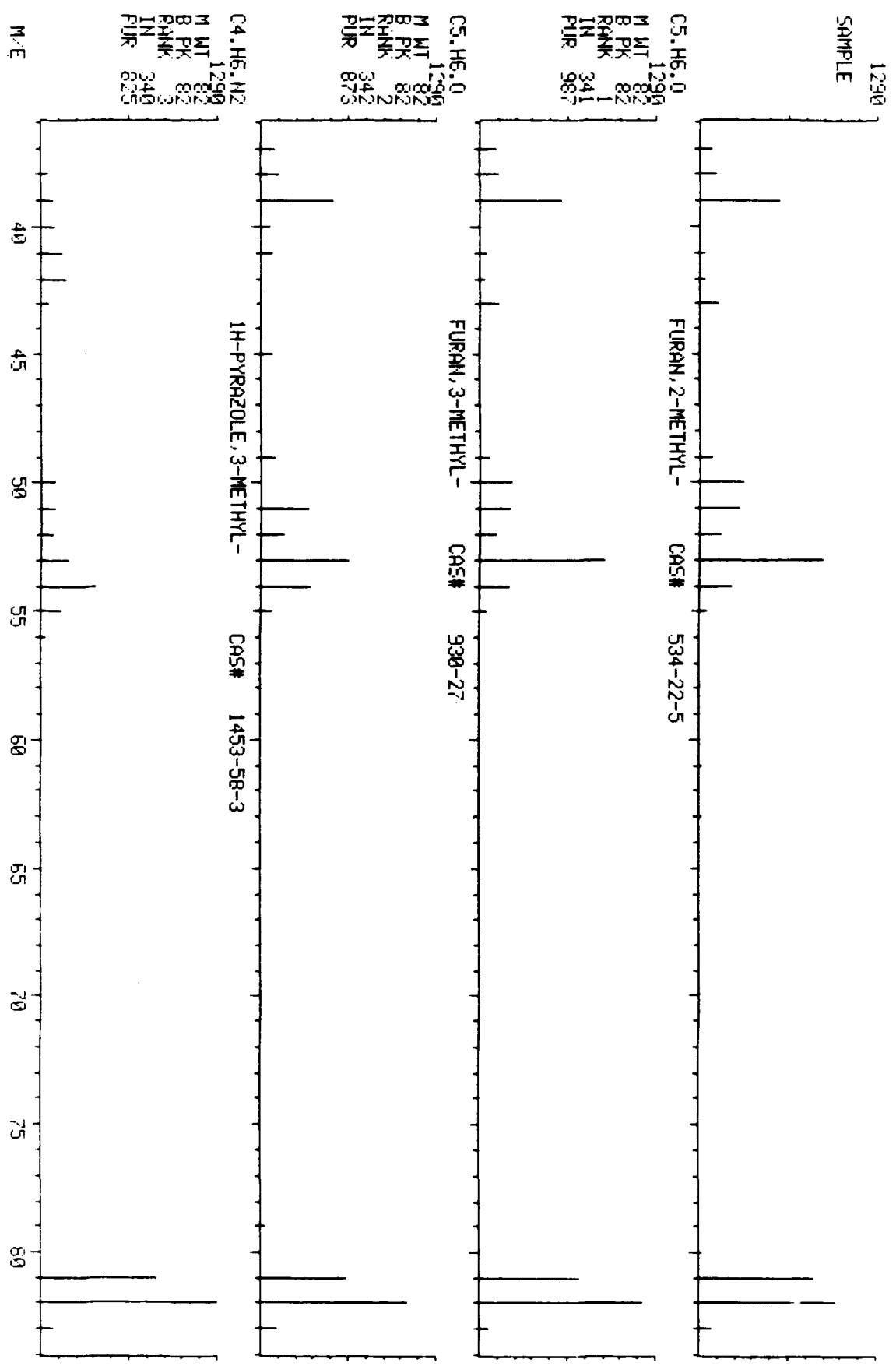


LIBRARY SEARCH
 12/27/89 7:10:00 + 4:58
 SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR ON#19
 ENHANCED (S 158 2N 0T)

COMPUCHEN LABS

DATA: GR009687019 # 397

BASE M/E: 82
 RIC: 508415.



LIBRARY SEARCH
12/27/89 7:10:00 + 2:14
SAMPLE: 1G CF#309687 CASE#18756.7 EPA#B202TAR ON#19
ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

DATA: GR009687C19 # 659

BASE M/E: 95
PIC: 610303.

1154
SAMPLE

C6.H8.O
1154
M WT 96
3 PK 96
RANK 776
IN 1
PUR 912

FURAN, 2,5-DIMETHYL- CAS# 625-86-5

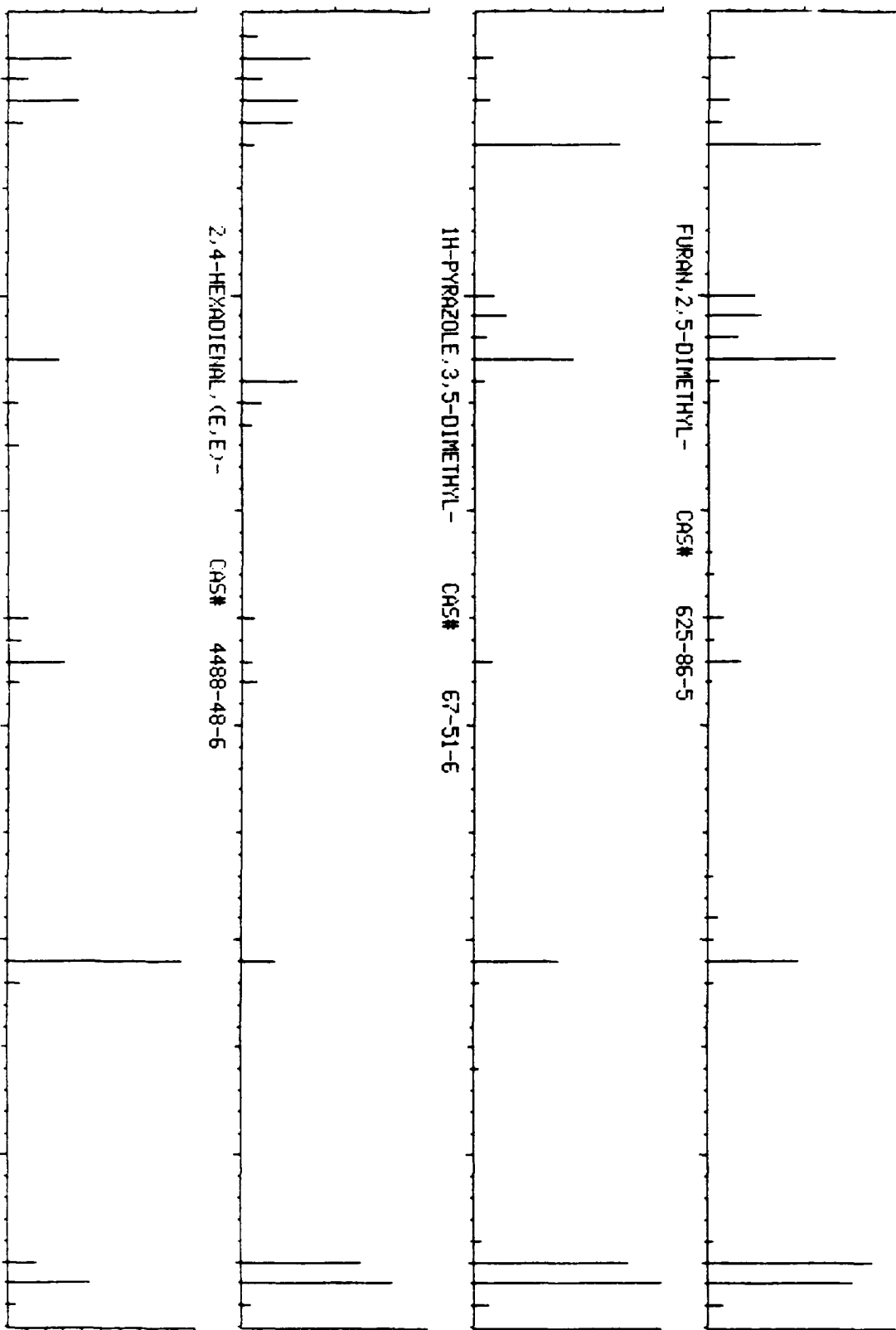
C5.H8.N2
1154
M WT 96
8 PK 96
RANK 769
IN 2
PUR 610

1H-PYRAZOLE, 3,5-DIMETHYL- CAS# 67-51-6

C6.H8.O
1154
M WT 96
8 PK 81
RANK 781
IN 3
PUR 606

2,4-HEXADIENAL, (E,E)- CAS# 4488-48-6

M/E 40 50 60 70 80 90



LIBRARY SEARCH
 12/27/89 7:10:00 + 11:03
 SAMPLE: 16 CC#309687 CASE#18756.7 EPA#B202TAR ON#19
 ENHANCED (5 156 2N 0T)

COMPUCHER LABS

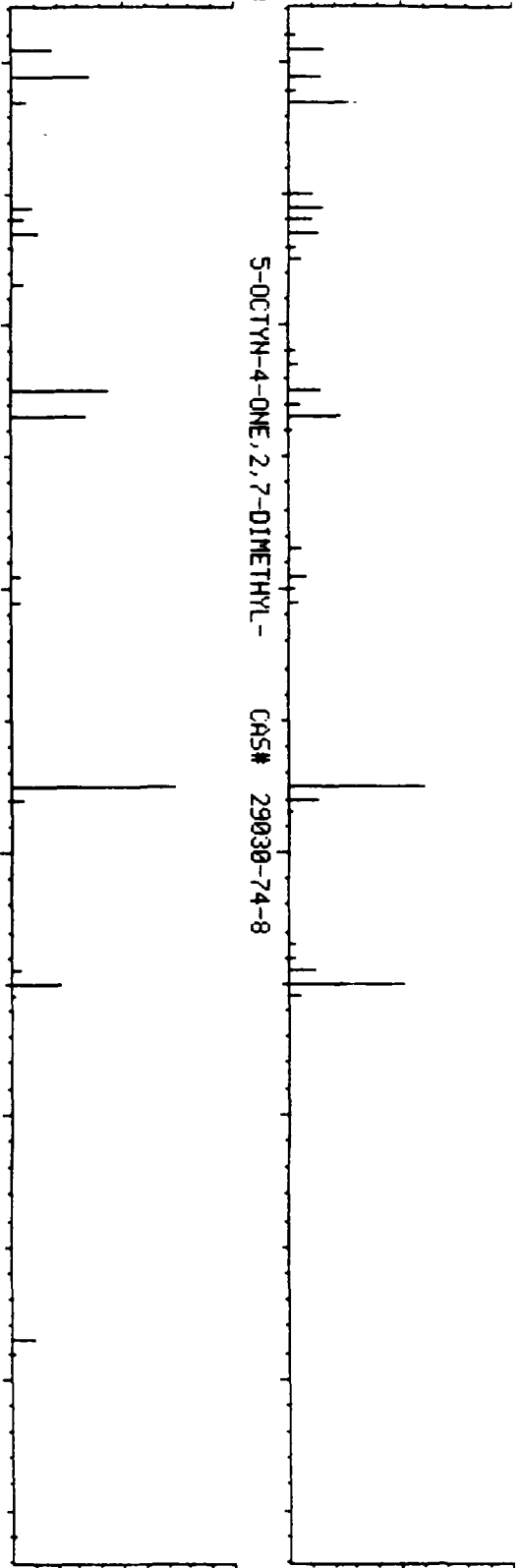
DATA: GR009687C19 # 884

BASE M/E: 95
 RID: 594943.

1659
 SAMPLE

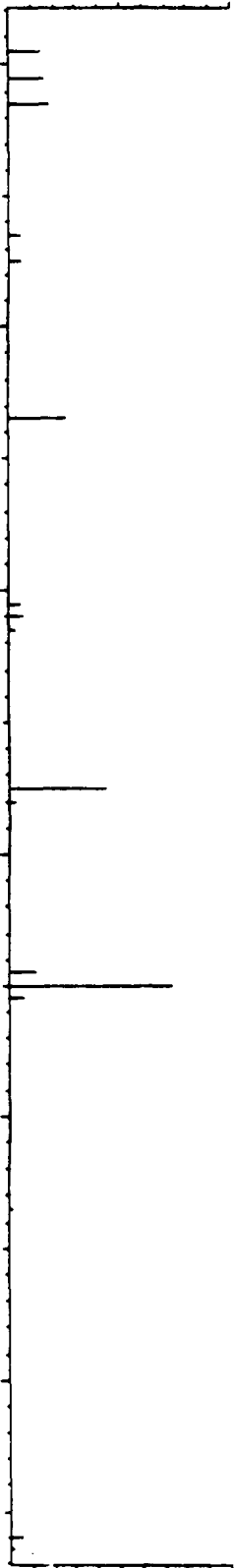
C10.H16.0
 M WT 1659
 B PK 95
 RANK 1
 IN 6245
 PUR 706

5-OCTYN-4-ONE, 2,7-DIMETHYL- CAS# 29030-74-8



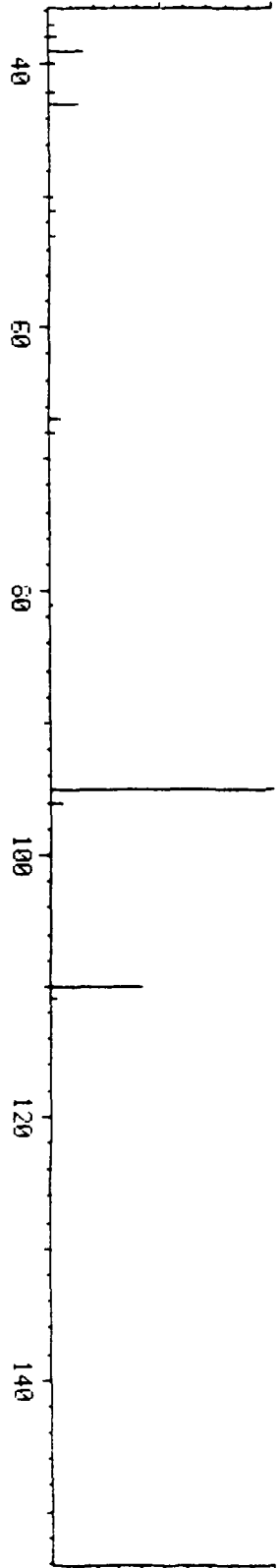
C9.H14.03
 M WT 1559
 B PK 110
 RANK 2
 IN 9072
 PUR 671

2-HEXENOICACID, 3,4,4-TRIMETHYL-5-OXO-, (Z)- CAS# 14919-56-3



C6.H6.O2
 M WT 1659
 B PK 95
 RANK 3
 IN 1596
 PUR 654

ETHANONE, 1-(2-FURANYL)- CAS# 1192-62-7

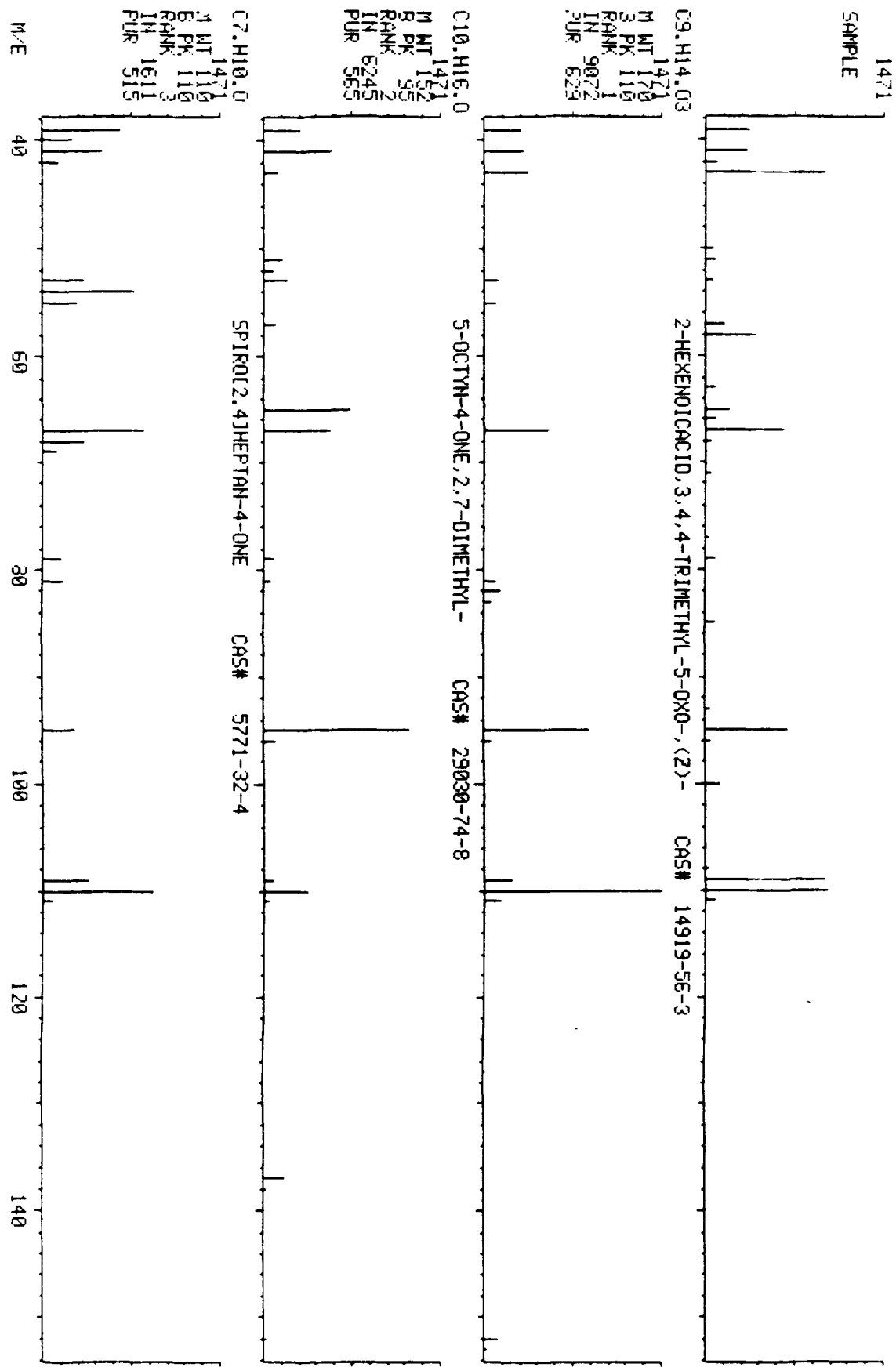


LIBRARY SEARCH
 12/27/89 7:10:00 + 11:30
 SAMPLE: 16 CC#309687 CASE#18756.7 EPAB202TAR ON#19
 ENHANCED (5 158 2H 0T)

COMPUchem LABS

DATA: GR009687019 # 920

BASE M/E: 110
 RIC: 560127.

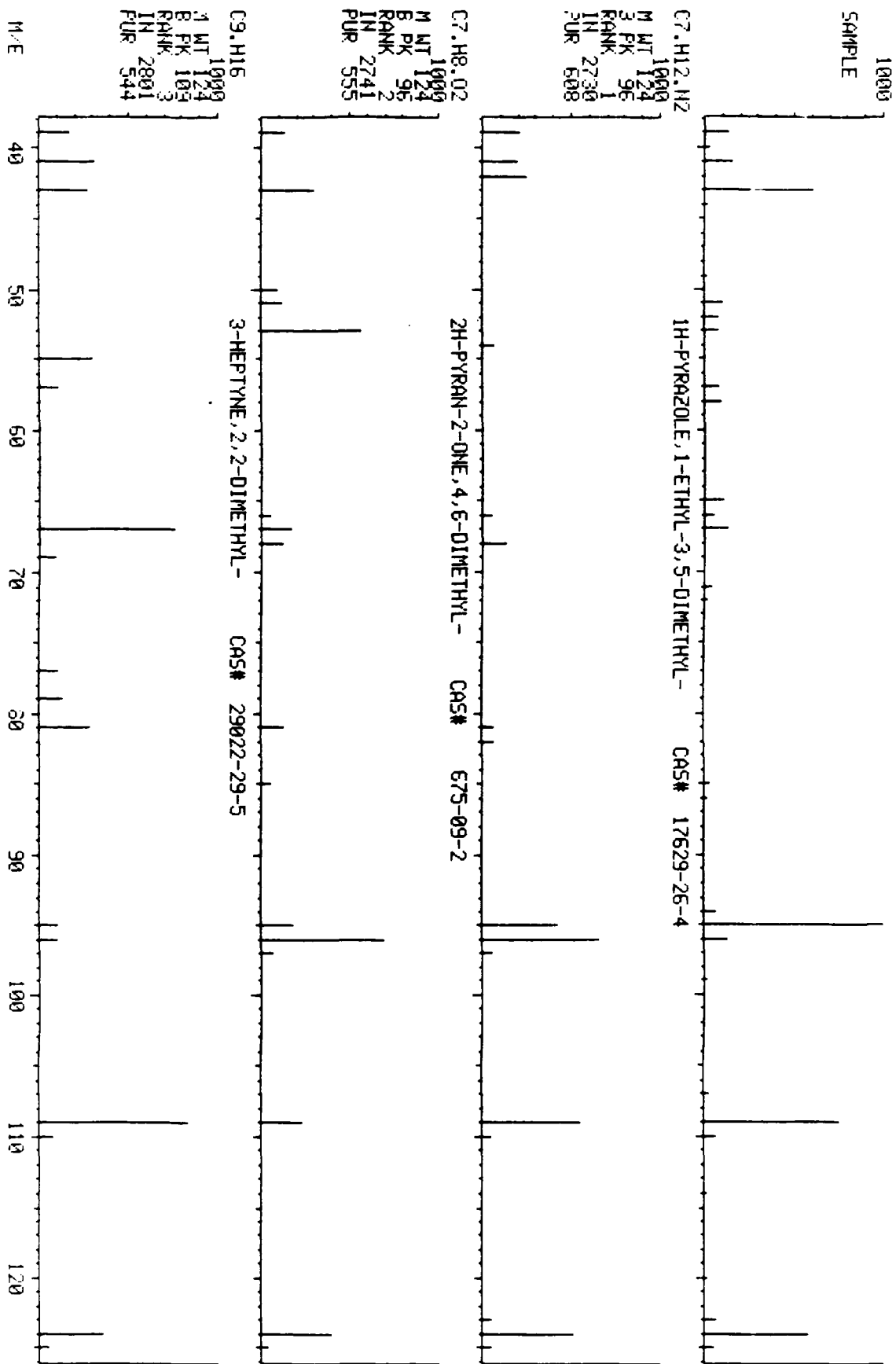


LIBRARY SEARCH
12/27/89 7:10:00 + 13:29
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR ON#19
ENHANCED (5 158 2N 0T)

COMPUchem LABS

DATA: GR009687C19 #1079

BASE M/E: 95
RIC: 376831.

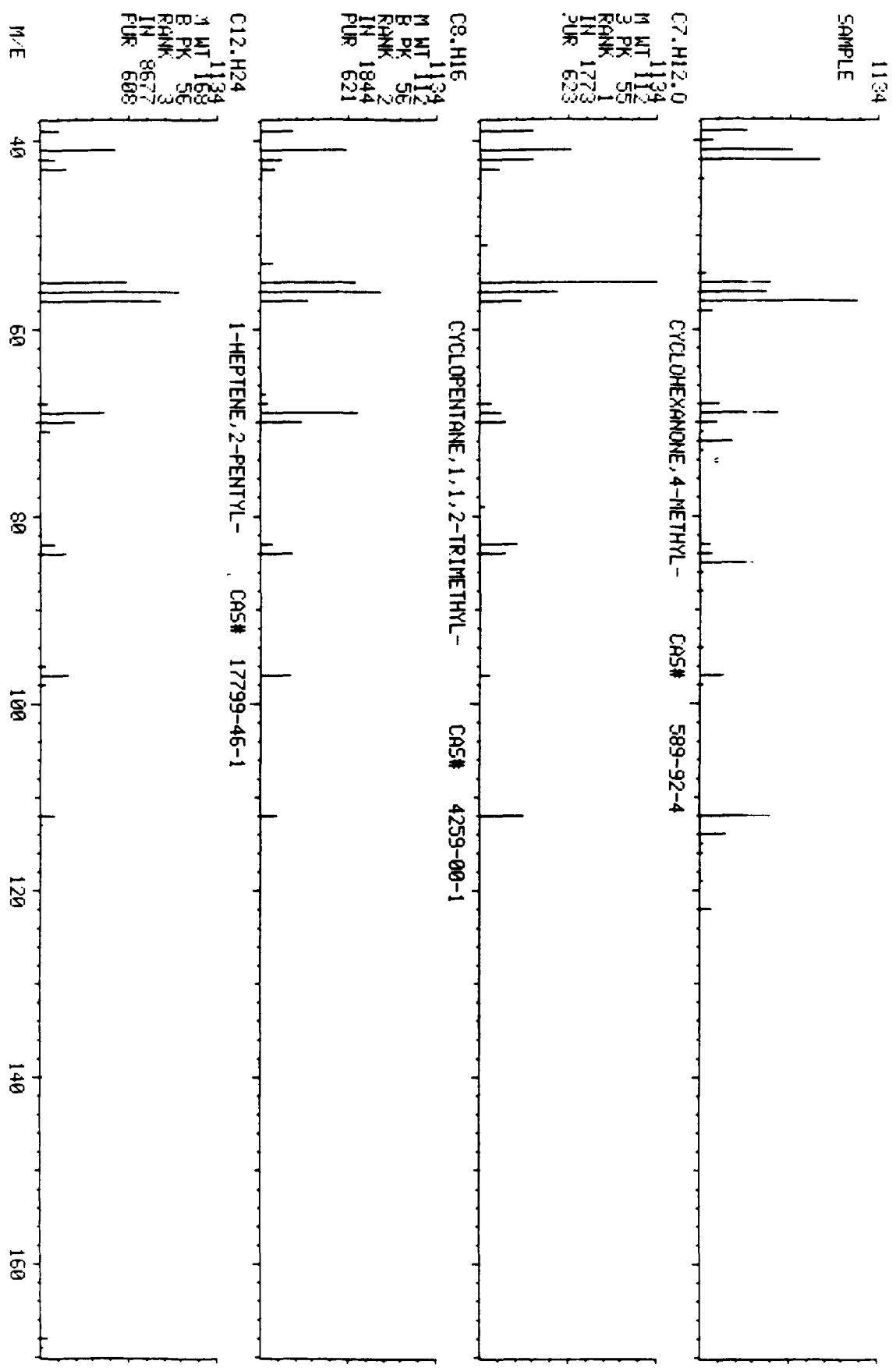


LIBRARY SEARCH
 12/27/89 7:10:00 + 14:05
 SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B202TAR QN#19
 ENHANCED (5 158 21 01)

COMPUchem LABS

DATA: GR009687C15 #1127

BASE M/E: 57
 RIC: 273407.

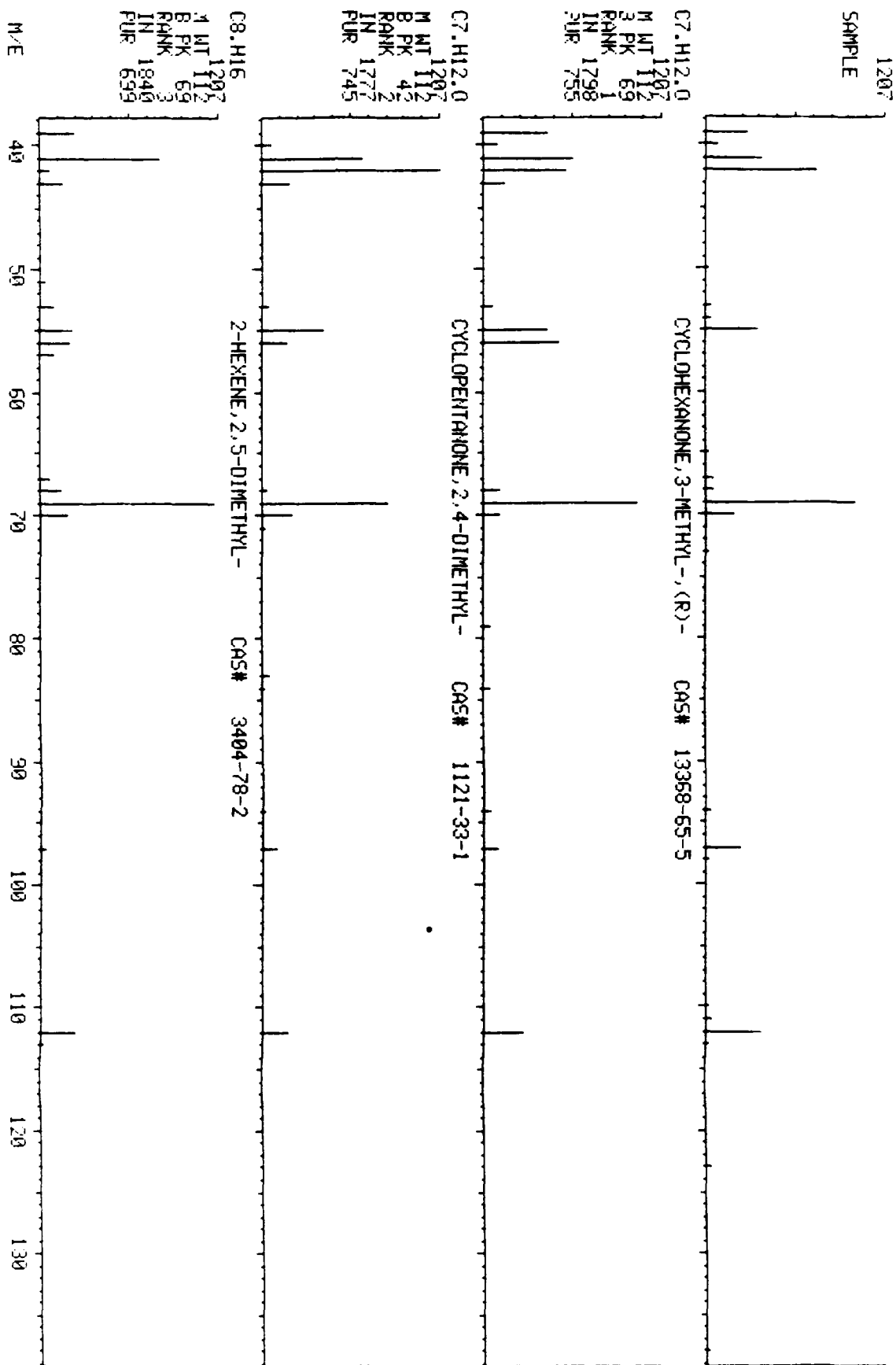


LIBRARY SEARCH
12/27/89 7:10:00 + 14:28
SAMPLE: 1G CC#309687 CASE#18756.7 EPA#B2021AR QM#19
ENHANCED (S 158 2H 0T)

COMPUCHEN LABS

DATA: GR09687019 #1157

BASE M/E: 69
RID: 285185.

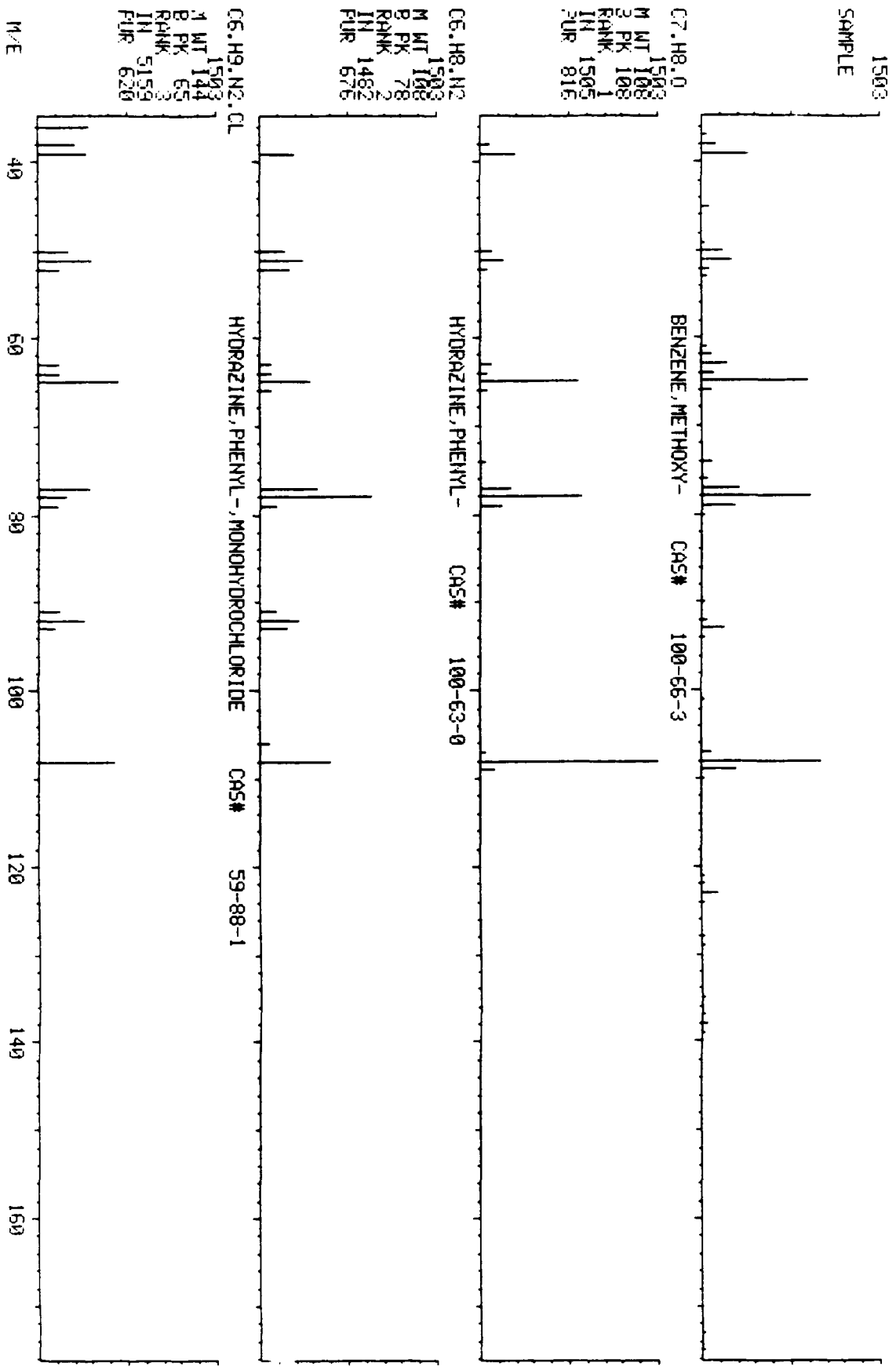


LIBRARY SEARCH
 12/27/89 7:16:00 + 14:45
 SAMPLE: 16 CC#809687 CASE#18756.7 EPA#B202TAR QN#19
 ENHANCED (S 156 2H 0T)

COMPUCHEN LABS

DATA: GR003687013 #1181

BASE M/E: 108
 RIC: 686079.



LIBRARY SEARCH
12/27/89 7:10:00 + 15:40
SAMPLE: 16-00#309687 CASE#18756.7 EPA#B202TAR Q#19
ENHANCED (5 158 2N 0T)

COMPUCHEN LABS

DATA: GR009687019 #1253

BASE M/E: 122
R/C: 48375.

1722
SAMPLE

C9.H12.02

M WT 1722
R PK 43
RANK 1
IN 6367
PUR 562

HYDROPEROXIDE, 1-METHYL-1-PHENYLETHYL

CAS#

80-15-9

C9.H12

M WT 1722
R PK 105
RANK 2
IN 2539
PUR 488

2,3-HEPTADIEN-5-YNE, 2,4-DIMETHYL-

CAS# 41898-89-9

C9.H12

M WT 1722
R PK 105
RANK 3
IN 2539
PUR 469

1,3-CYCLOPENTADIENE, 5-(1-METHYLPROPYLIDENE)-

CAS# 3141-02-4

M/E 40 50 60 70 80 90 100 110 120 130

LIBRARY SEARCH
 12/27/89 7:10:00 + 16:06
 SAMPLE: 1G CC#309687 CASE#18756.7 EPA#E2021AR QM#19
 ENHANCED (S 158 ZN 01)

COMPUJEN LABS

DATA: GR009687C19 #1288

BASE M/E: 105
 RID: 469503.

1605
 SAMPLE

C9.H12
 1605

M WT 120
 B PK 105
 RANK 1
 IN 2535
 PUR 787

1,3-CYCLOPENTADIENE, 5-(1-METHYLPROPYLIDENE) - CAS# 3141-02-4

C9.H12
 1605

M WT 120
 B PK 105
 RANK 2
 IN 2523
 PUR 784

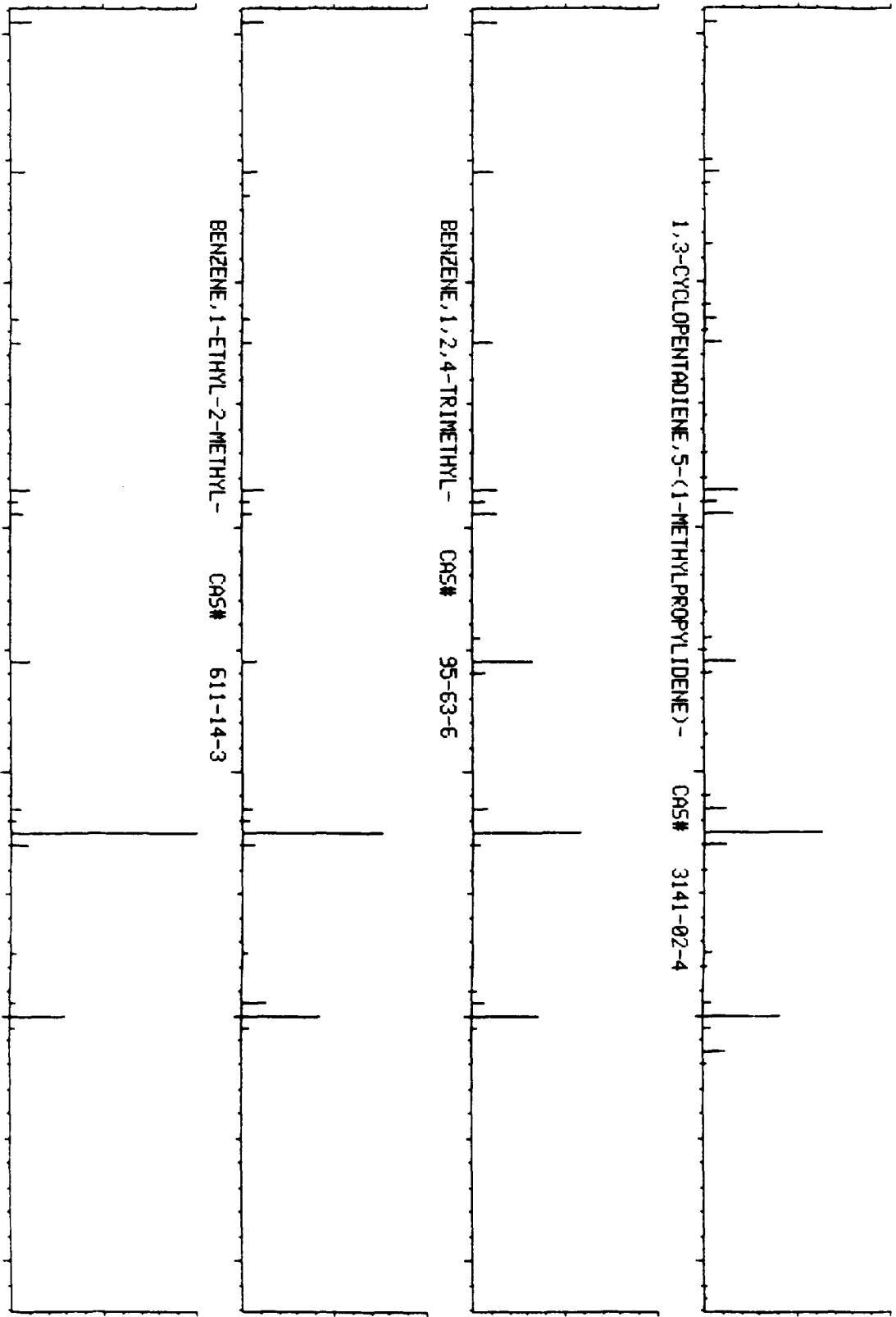
BENZENE, 1,2,4-TRIMETHYL - CAS# 95-63-6

C9.H12
 1605

M WT 120
 B PK 105
 RANK 3
 IN 2528
 PUR 784

BENZENE, 1-ETHYL-2-METHYL - CAS# 611-14-3

M/E 40 50 80 100 120 140



LAB INSTRUCTIONS

SAMPLED DATE 12/18/89

RECEIPT DATE 12/20/89

CASE# 18756 7

DUE DATE

VQA
GC/MS WORKSHEET

COMPUCHEM# 309687R

REX] R2C] DC] ()
R3C] R4C] DEC] ()

L L SOLID, EPA SOW 2/88

Sample Prep Code---155
Instrument Code---417
Compound List-----494
Surrogate Std-----394
Internal Std-----036

=====

SAMPLE ID# B202TAR Dry Wt. Factor 2.78 % Moisture 64

=====

GC/MS ANALYSIS

Amount Purged: [] 10mls/Xg soil or [] Dilution _____ ul/10000ul/Xg soil
Internal Standard Volume Added 5 ul
Surrogate Standard Volume Added 5 ul
BFB Filename BF891227C19 Disk (2858A)
Blank Filename 614010660019 Disk ()
Standard Filename 67891227C19 Disk (8)
Sample Filename 6A024687C19 Disk ()

ANALYST(S) Injection 1977-16 Work-up 1172-16

GC/MS REVIEW

CONDITION
CODE

DW
DI

Entry Codes OK, EA, ES, SM, JS, SL, SH, JA, DA

Non-Entry Codes IM, IL, IH, SW, CT, CS, PC, NR
IF, LA, DI, CO, RN, DW, SI, SF
UP, BB, OT, VC, FO, NS

Disposition: [] Complete
[] Reprep neat required
[] Reprep using _____
[] Dilute ()
 Reprep neat

Extraneous Peak Search Results:
of Peaks Found: 10

Quality Assurance Notice(s):
Notices Required _____

COMMENTS:

GC/MS Review [Signature] Date 1/3/90 Auditor _____ Date _____

REPORT INTEGRATION
Final Reportable Package(s): 6R009687C19 / C3R09687C13 Total # of Injections _____

QA COMMENTS:

=====

INITIALS _____ DATE _____

=====

FINAL REVIEW INITIALS _____ DATE _____

AC1007 (05/89)

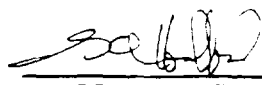
VOLATILE PREPARATION WORKSHEET

QUEUE # 19 PREP CODE 155 ASSIGNED TO Hanshad Joshi DATE 12/22/89


SAMPLE NUMBER	CASE NUMBER	SDG	QC SAMPLE		SAMPLE WEIGHT (G) VOLUME (ML)	SAMPLE ID	COMMENTS
			TYPE	ORIGINAL			
309679R	18756	0007			1.0g	B201A	
309680R			SS	309679	1.0g	B201A MS	
309681R			SS	309679	1.0g	B201A MSD	
309682R			BS		0.0g		
309686R					1.0g	B201B	
309687R					1.0g	B202AR	
310543	18244	0142			5.0g	MW-17-30	
310555	18477	0132			5.0g	EPTP10122	
310560			SS	310555	5.0g	EPTP10122 MS	
310561			SS	310555	5.0g	EPTP10122 MSD	
310562			BS		0.0g		
310659			B1		5.0ml	F4	
310660			B2		0.0ml	F5	
310661			B3		0.0ml	F6	
310662			B4		0.0ml	F7	
310663			B5		0.0ml	F8	

SURROGATE # LOT # MANUAL OPERATOR 73, 518
 AMOUNT
 RELINQUISHED BY DATE 12/22/89 RECEIVED BY 182 N/A DATE

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	463	68100	50.0		
221	50	CHLOROMETHANE				BDL	
231	62	VINYL CHLORIDE				BDL	
220	94	BROMOMETHANE				BDL	
209	64	CHLOROETHANE				BDL	
216	96	1,1-DICHLOROETHENE				BDL	
254	76	CARBON DISULFIDE				BDL	
252	43	ACETONE (2-PROPANONE)			250.0	³⁵⁰⁰ 1200	E 13
248	114 I	1,4-DIFLUOROBENZENE (IS)	608	245000	50.0		
222	84	METHYLENE CHLORIDE			599.0	⁸³⁰⁰ 3000	E 13
226	96	TRANS-1,2-DICHLOROETHENE				BDL	
214	63	1,1-DICHLOROETHANE				BDL	
257	43	VINYL ACETATE				BDL	
237	96	CIS-1,2-DICHLOROETHENE				BDL	
253	72	2-BUTANONE			362.0	⁵⁰⁰⁰ 1800	E
211	83	CHLOROFORM				BDL	
227	97	1,1,1-TRICHLOROETHANE				BDL	
206	117	CARBON TETRACHLORIDE				BDL	
203	78	BENZENE			88.1	¹²⁰⁰ 440	
215	62	1,2-DICHLOROETHANE				BDL	
270	117 I	D5-CHLOROBENZENE (IS) RO#29	1002	246000	50.0		
229	130	TRICHLOROETHENE				BDL	
217	63	1,2-DICHLOROPROPANE			17.2	88 BDL	
212	83	BROMODICHLOROMETHANE				BDL	
218	75	CIS-1,3-DICHLOROPROPENE				BDL	
256	43	4-METHYL-2-PENTANONE			86.9	¹²⁰⁰ 430	
225	92	TOLUENE			622.0	⁸⁶⁰⁰ 3100	E
250	75	TRANS-1,3-DICHLOROPROPENE			8.2	400 BDL	
228	97	1,1,2-TRICHLOROETHANE				BDL	
224	164	TETRACHLOROETHENE				BDL	
255	43	2-HEXANONE			838.0	¹²⁰⁰⁰ 4200	E 13
208	129	DIBROMOCHLOROMETHANE				BDL	
207	112	CHLOROBENZENE				BDL	
219	106	ETHYLBENZENE			575.0	⁸⁰⁰⁰ 2900	E
330	106	M, P-XYLENE			532.0	⁷⁴⁰⁰ 2700	E
239	106	O-XYLENE			568.0	⁷⁴⁰⁰ 2800	E
251	104	STYRENE			164.0	²³⁰⁰ 820	
205	173	BROMOFORM				BDL	
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	
258	65 S	D4-1,2-DICHLOROETHANE RO#57			48.5	97 %	
247	95 S	BROMOFLUOROBENZENE			46.7	93 %	
233	98 S	D8-TOLUENE RO#59			44.1	88 %	
289	106	XYLENES (TOTAL)			1100.0	5500 ¹⁵⁰⁰⁰	E

CORRECTED/REVIEWED BY 
(GC/MS DATA REVIEWER)
DATE 12/31/95

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
299	96	1,2-DICHLOROETHENE (TOTAL)				BDL	2
CHECKSUMS:							
		3979.	2073	559100.	6099.7	29295	

CORRECTED/REVIEWED BY 
(GC/MS DATA REVIEWER)

DATE 12/3/57

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P
40	258	D4-1,2-DICHLOROETHANE RO#57	48.5	50.0	97.	70-121	X
41	247	BROMOFLUOROBENZENE	46.7	50.0	93.	74-121	X
42	233	D8-TOLUENE RO#59	44.1	50.0	88.	81-117	X

* ADVISORY SURROGATE ONLY

++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

CORRECTION FACTOR CALCULATION:

$$\frac{5.0 \text{ G}}{\text{WET WEIGHT OF SAMPLE (G)}} \times \frac{\text{GC/MS DILUTION FACTOR}}{\text{GC/MS DILUTION FACTOR}} \times \text{DRY WEIGHT FACTOR} =$$

$$\frac{5.0 \text{ G}}{1.00 \text{ (G)}} \times \frac{1.0}{1.00} \times \frac{2.78}{1.00} = 13.9$$

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:

THE SURROGATES ARE ADDED TO THE SAMPLE PRIOR TO SPARGING.
SURROGATE SPIKE CONVERSION FACTOR = 1.

VERSION 8

CORRECTED/REVIEWED BY *Saltzman*
(GC/MS DATA REVIEWER)

DATE 12-31-

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

B202TARDL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C3R09687C13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 Date Analyzed: 12/29/89
 Column: (pack/cap) CAP Dilution Factor: 3.3

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>	Q
74-87-3	-----Chloromethane	11000	U
74-83-9	-----Bromomethane	11000	U
75-01-4	-----Vinyl Chloride	11000	U
75-00-3	-----Chloroethane	11000	U
75-09-2	-----Methylene Chloride	16000	BD
67-64-1	-----Acetone	15000	D
75-15-0	-----Carbon Disulfide	5700	U
75-35-4	-----1,1-Dichloroethene	5700	U
75-34-3	-----1,1-Dichloroethane	5700	U
540-59-0	-----1,2-Dichloroethene (total)	5700	U
67-66-3	-----Chloroform	5700	U
107-06-2	-----1,2-Dichloroethane	5700	U
78-93-3	-----2-Butanone	11000	U
71-55-6	-----1,1,1-Trichloroethane	5700	U
56-23-5	-----Carbon Tetrachloride	5700	U
108-05-4	-----Vinyl Acetate	11000	U
75-27-4	-----Bromodichloromethane	5700	U
78-87-5	-----1,2-Dichloropropane	5700	U
10061-01-5	-----cis-1,3-Dichloropropene	5700	U
79-01-6	-----Trichloroethene	5700	U
124-48-1	-----Dibromochloromethane	5700	U
79-00-5	-----1,1,2-Trichloroethane	5700	U
71-43-2	-----Benzene	5000	DJ
10061-02-6	-----Trans-1,3-Dichloropropene	5700	U
75-25-2	-----Bromoform	5700	U
108-10-1	-----4-Methyl-2-Pentanone	11000	U
591-78-6	-----2-Hexanone	11000	U
127-18-4	-----Tetrachloroethene	5700	U
79-34-5	-----1,1,2,2-Tetrachloroethane	5700	U
108-88-3	-----Toluene	100000	D
108-90-7	-----Chlorobenzene	5700	U
100-41-4	-----Ethylbenzene	100000	D
100-42-5	-----Styrene	22000	D
1330-20-7	-----Total Xylenes	280000	DE

FORM I VOA

1/87 Rev.

1E
VOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

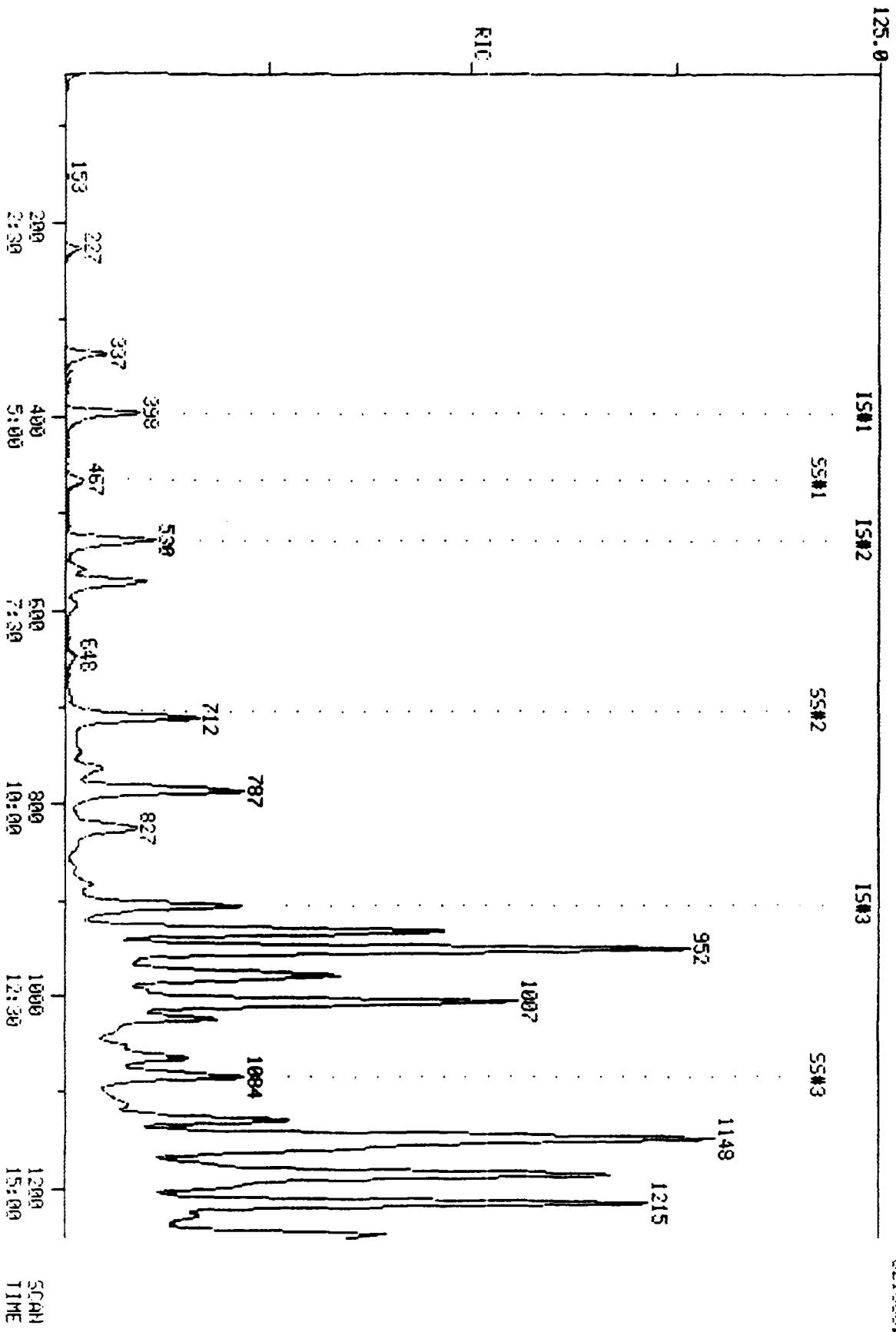
B202TARDL

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07
 Matrix: (soil/water) SOIL Lab Sample ID: 309687
 Sample wt/vol: 4.0 (g/mL) G Lab File ID: C3R09687C13
 Level: (low/med) MED Date Received: 12/20/89
 % Moisture: not dec. 64 Date Analyzed: 12/29/89
 Column (pack/cap) CAP Dilution Factor: 3.3

Number TICs found: 10 CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 534-22-5	FURAN, 2-METHYL-	4.20	29000	J
2. 625-86-5	FURAN, 2, 5-DIMETHYL-	7.15	49000	J
3. 1192-62-7	ETHANONE, 1-(2-FURANYL)-	9.84	61000	J
4. 37609-41-9	BICYCLO[3.2.1]OCT-2-ENE, 3-(1	12.24	85000	J
5. 24156-95-4	2-CYCLOPENTEN-1-ONE, 3, 5, 5-TR	12.80	40000	J
6.	UNKNOWN	13.30	34000	J
7.	SUBST. BENZENE	14.10	48000	J
8.	SUBST. BENZENE	14.35	300000	J
9.	SUBST. BENZENE	14.82	220000	J
10.	SUBST. BENZENE	15.19	170000	J

RIC
 12/29/89 1:52:00
 SAMPLE: 30UL CC#309687 EPA#B2021AR RE CASE#18756 ON #13
 COND5.:
 COMPUTHEM LABS
 DL SA# 112190
 COMPUTHEM DATA: C3R09687013 SCANS 47 TO 1250



QUANTITATION REPORT FILE: C3R09687C13
 DATA: C3R09687C13.TI
 12/29/89 1:52:00 *DL sent 1/20/90*
 SAMPLE: 30UL CC#309687 EPA#B202TAR RE CASE#18756 DN #13
 CONDS.:
 SUBMITTED BY: 13 ANALYST: 1452

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
 RESP. FAC. FROM LIBRARY ENTRY

- NO NAME
- 1 *234 BROMOCHLOROMETHANE (IS) <75-97-5> NA#1
- 2 221 CHLOROMETHANE <74-87-3> NA#2
- 3 231 VINYL CHLORIDE <75-01-4> NA#3
- 4 220 BROMOMETHANE <78-83-9> NA#4
- 5 209 CHLOROETHANE <75-00-3> NA#5
- 6 216 1,1-DICHLOROETHENE <75-35-4> NA#8
- 7 254 CARBON DISULFIDE <75-15-0> NA#9
- 8 252 ACETONE (2-PROPANONE) <67-64-1> NA#13
- 9 *248 1,4-DIFLUOROBENZENE (IS) <340-36-3> NA#14
- 10 222 METHYLENE CHLORIDE <75-09-2> NA#16
- 11 226 TRANS-1,2-DICHLOROETHENE <156-60-5> NA#17
- 12 214 1,1-DICHLOROETHANE <75-34-3> NA#19
- 13 257 VINYL ACETATE <108-05-4> NA#20
- 14 237 CIS-1,2-DICHLOROETHENE <156-59-2> NA#21
- 15 253 2-BUTANONE <78-93-3> NA#22
- 16 211 CHLOROFORM <67-66-2> NA#23
- 17 227 1,1,1-TRICHLOROETHANE <71-55-6> NA#24
- 18 206 CARBON TETRACHLORIDE <56-23-5> NA#25
- 19 203 BENZENE <71-43-2> NA#26
- 20 215 1,2-DICHLOROETHANE <107-06-2> NA#27
- 21 *270 D5-CHLORO BENZENE (IS) <XX-XX-X> NA#29
- 22 229 TRICHLOROETHENE <79-01-6> NA#30
- 23 217 1,2-DICHLOROPROPANE <78-87-5> NA#31
- 24 212 BROMODICHLOROMETHANE <75-27-4> NA#33
- 25 218 CIS-1,3-DICHLOROPROPENE <10061-1-5> NA#35
- 26 256 4-METHYL-2-PENTANONE <108-01-1> NA#36
- 27 225 TOLUENE <108-88-3> NA#37
- 28 250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> NA#38
- 29 228 1,1,2-TRICHLOROETHANE <79-00-5> NA#39
- 30 224 TETRACHLOROETHENE <127-18-4> NA#41
- 31 255 2-HEXANONE <591-78-6> NA#42
- 32 208 DIBROMOCHLOROMETHANE <124-48-1> NA#43
- 33 207 CHLOROBENZENE <108-90-7> NA#45
- 34 219 ETHYLBENZENE <100-41-4> NA#47
- 35 330 M,P-XYLENE <133-02-7> NA#48
- 36 239 O-XYLENE <133-02-7> NA#49
- 37 251 STYRENE <100-42-5> NA#50
- 38 205 BROMOFORM <75-25-2> NA#51
- 39 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> NA#54
- 40 *258 D4-1,2-DICHLOROETHANE NA#57
- 41 *247 BROMOFLUOROBENZENE <460-00-4> NA#58
- 42 *233 D8-TOLUENE NA#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	397	4:58	1	1.000	A BB	49580.	50.000 UG/L	7.50
2	50	NOT FOUND							

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
3	62	NOT FOUND							
4	94	NOT FOUND							
5	64	NOT FOUND							
6	96	NOT FOUND							
7	76	NOT FOUND							
8	43	NOT FOUND	195						
9	114	530	6:37	9	1.000	A BB	172864. 3369	50.000 UG/L 13,220 µg/L	7.50
10	84	227	2:50	1	0.572	A BV	20440.	14.102 UG/L	2.11
11	96	NOT FOUND							
12	63	NOT FOUND							
13	43	NOT FOUND							
14	96	NOT FOUND							
15	72	NOT FOUND							
16	83	NOT FOUND							
17	97	NOT FOUND							
18	117	NOT FOUND							
19	78	464	5:48	9	0.875	A BB	13235.	4.329 UG/L	0.65
20	62	NOT FOUND							
21	117	906	11:19	21	1.000	A BB	250054.	50.000 UG/L	7.50
22	130	NOT FOUND							
23	63	NOT FOUND							
24	83	NOT FOUND							
25	75	NOT FOUND							
26	43	714	8:55	21	0.788	A*BB	2827.	2.424 UG/L	0.36
27	92	712	8:54	21	0.786	A BB	151864.	91.345 UG/L	13.70
28	75	NOT FOUND							
29	97	786	9:49	9	1.483	A BB	1846.	1.930 UG/L	0.29
30	164	NOT FOUND							
31	43	NOT FOUND							
32	129	NOT FOUND							
33	112	NOT FOUND							
34	106	933	11:40	21	1.030	A BV	186205.	90.394 UG/L	13.55
35	106	952	11:54	21	1.051	A VB	526501.	148.536 UG/L	22.27
36	106	1006	12:34	21	1.110	A BB	306746.	98.871 UG/L	14.82
37	104	1012	12:39	21	1.117	A BB	105930.	19.536 UG/L	2.93
38	173	NOT FOUND							
39	83	NOT FOUND							
40	65	467	5:50	1	1.176	A BB	28655.	16.482 UG/L	2.47
41	95	1085	13:34	21	1.198	A VB	74996.	16.620 UG/L	2.49
42	98	704	8:48	21	0.777	A BB	33435.	12.410 UG/L	1.86

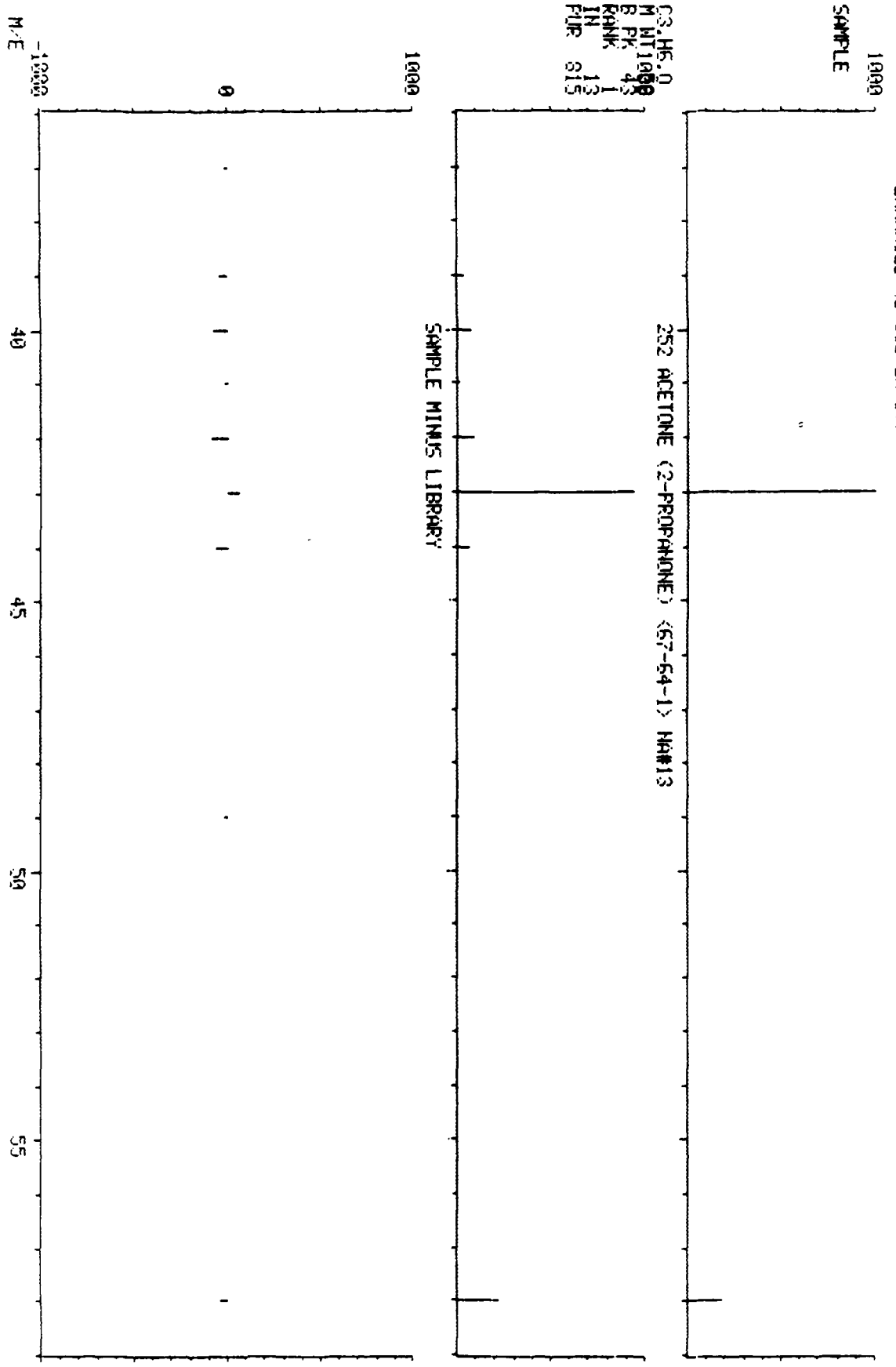
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:07	0.97	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	0:52		10.000			50.00		0.882	
3	0:58		10.000			50.00		1.005	
4	1:11		10.000			50.00		1.237	
5	1:18		10.000			50.00		0.626	
6	2:10		5.000			50.00		1.230	
7	2:16		5.000			50.00		3.765	
8	2:31		10.000			50.00		0.257	
9	6:46	0.98	5.000	0.20	50.00	50.00	1.000	1.000	1.00
10	2:58	0.95	5.000	0.11	14.10	50.00	0.412	1.462	0.28
11	3:18		5.000			50.00		1.270	
12	3:56		5.000			50.00		2.282	
13	4:14		10.000			50.00		0.848	
14	4:48		5.000			50.00		1.568	

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
15	5:01		10.000			50.00		0.124	
16	5:22		5.000			50.00		2.770	
17	5:24		5.000			50.00		0.713	
18	5:37		5.000			50.00		0.656	
19	5:57	0.97	5.000	0.18	4.33	50.00	0.077	0.884	0.09
20	6:06		5.000			50.00		1.804	
21	11:25	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
22	7:01		5.000			50.00		0.453	
23	7:22		5.000			50.00		0.350	
24	7:56		5.000			50.00		0.642	
25	8:40		5.000			50.00		0.553	
26	9:07	0.98	15.000	0.05	2.42	50.00	0.011	0.233	0.05
27	9:04	0.98	5.000	0.16	91.35	50.00	0.607	0.332	1.83
28	9:40		5.000			50.00		0.243	
29	9:56	0.99	5.000	0.30	1.93	50.00	0.011	0.277	0.04
30	9:53		5.000			50.00		0.303	
31	10:35		15.000			50.00		0.228	
32	10:30		5.000			50.00		0.846	
33	11:28		5.000			50.00		0.966	
34	11:49	0.99	5.000	0.21	90.39	50.00	0.745	0.412	1.81
35	12:04	0.99	5.000	0.21	148.54	50.00	2.106	0.709	2.97
36	12:46	0.99	5.000	0.22	98.87	50.00	1.227	0.620	1.98
37	12:50	0.99	5.000	0.22	19.54	50.00	0.424	1.084	0.39
38	13:06		5.000			50.00		0.762	
39	14:19		5.000			50.00		0.705	
40	6:00	0.97	5.000	0.24	16.48	50.00	0.578	1.753	0.33
41	13:44	0.99	5.000	0.24	16.62	50.00	0.300	0.902	0.33
42	8:57	0.98	5.000	0.16	12.41	50.00	0.134	0.539	0.25

LIBRARY SEARCH
12/29/89 1:52:00 + 2:26
SAMPLE: 3MUL 01#309687 EPA#B202TAR RE CASE#18756 ON #13
ENHANCED (S 158 2N 0T)

COMPUchem Labs
DATA: 03R09687013 # 195
BASE M/E: 43
R1C: 152

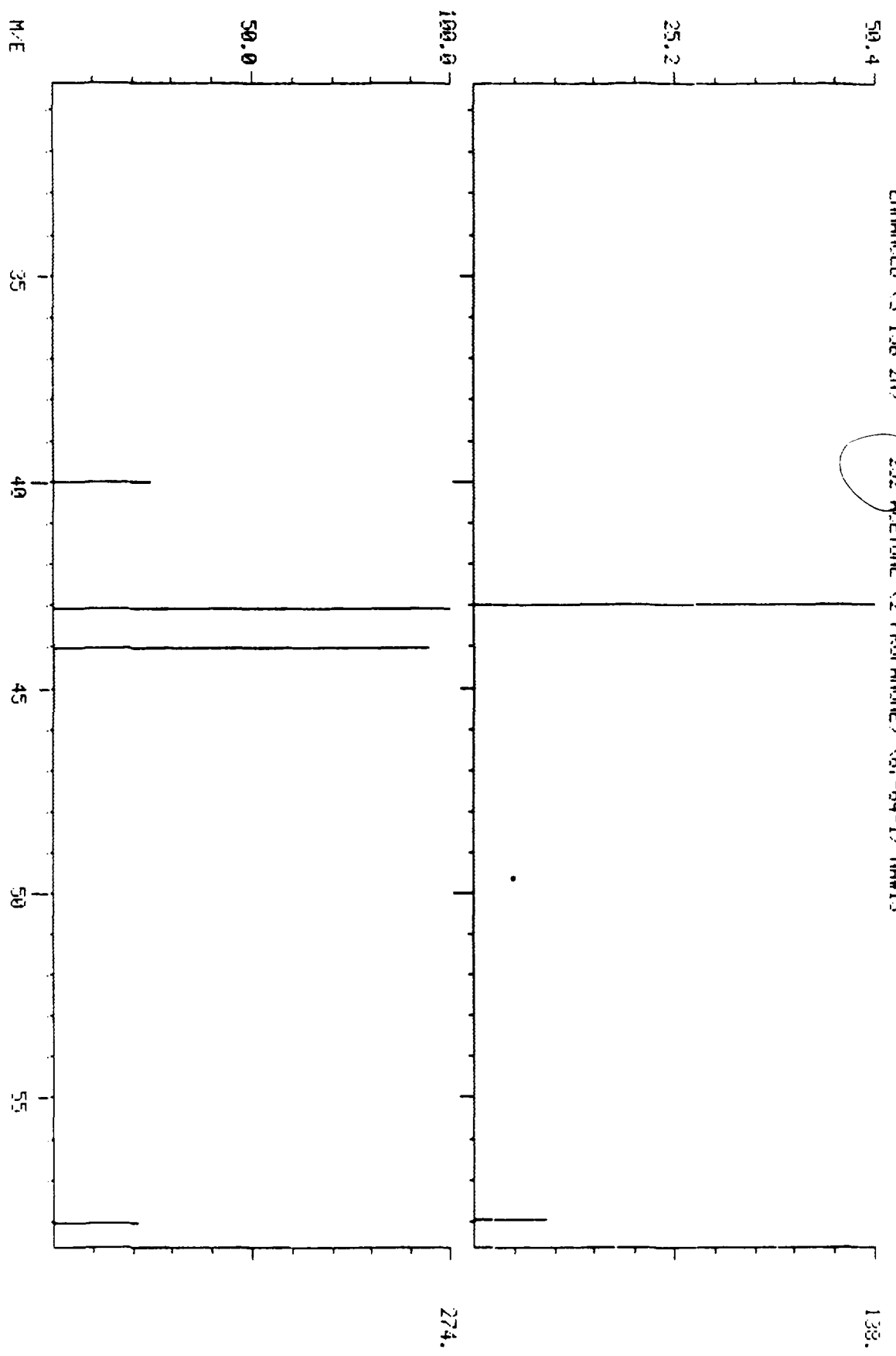
03.H5.0
M UT 1000
R PK 43
IN 13
FILR 815



DUAL MASS SPECTRUM
12/29/89 1:52:00 + 2:26
SAMPLE: 30UL CC#309687
ENHANCED (S 158 2H)

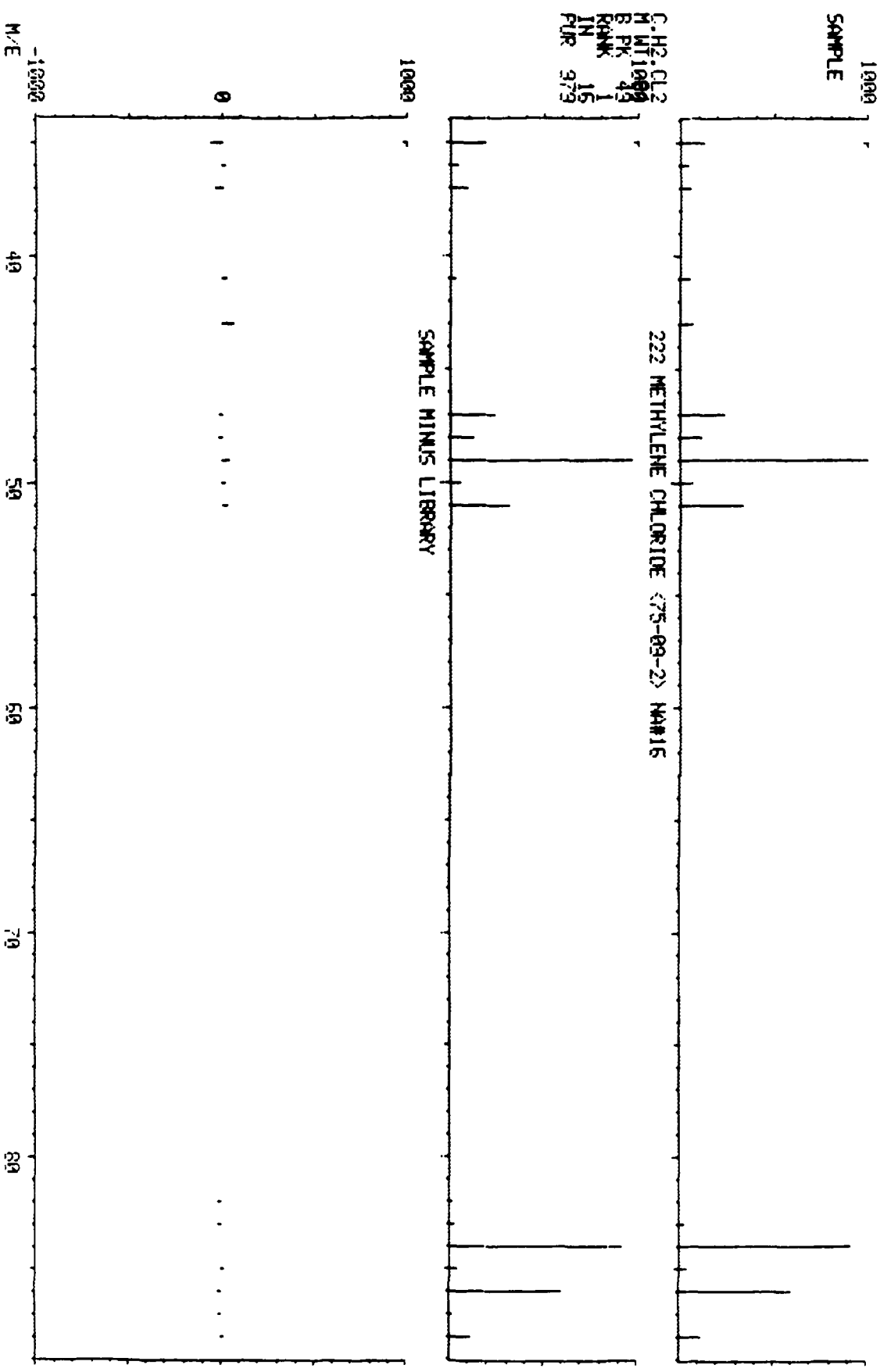
COMPUCHER LABS
DATA: C309687013 #195
252 ACETONE (2-PROPANONE) (57-54-1) NA#13

BASE M/E: 43
RIC: 162.7 657.



COMPUchem LABS
 LIBRARY SEARCH
 12/29/89 1:52:00 + 2:50
 SAMPLE: 30UL CC#309687 EPA#8202TAR RE CASE#18756 ON #13
 ENHANCED (S 158 2H 0T)
 DL 544.13/90
 DATA: C3R09687C13 # 227
 BASE M/E: 49
 RIC: 8415.

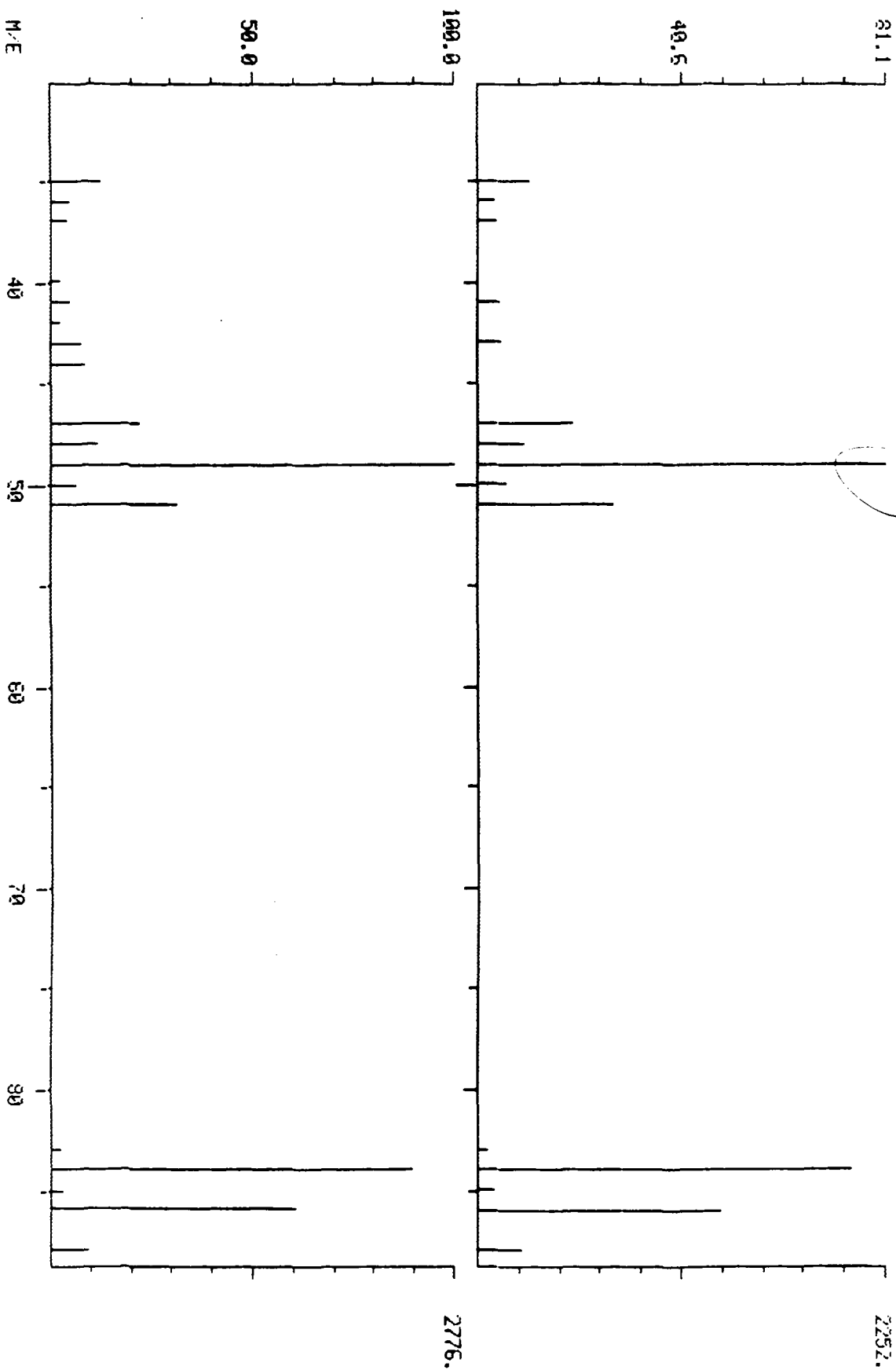
C.H2.C12
 H.M.I.1000
 R.P.K. 49
 R.H.A.K. 15
 T.N. 15
 P.U.R. 979



DUAL MASS SPECTRUM
12/29/89 1:52:00 + 2:50
SAMPLE: 30UL CC#309687
ENHANCED (5 158 2N)

COMPIGHEM LABS
DL *gaw lab*
DATA: C3R09687013 #227
222 METHYLENE CHLORIDE (75-09-2) NA#15

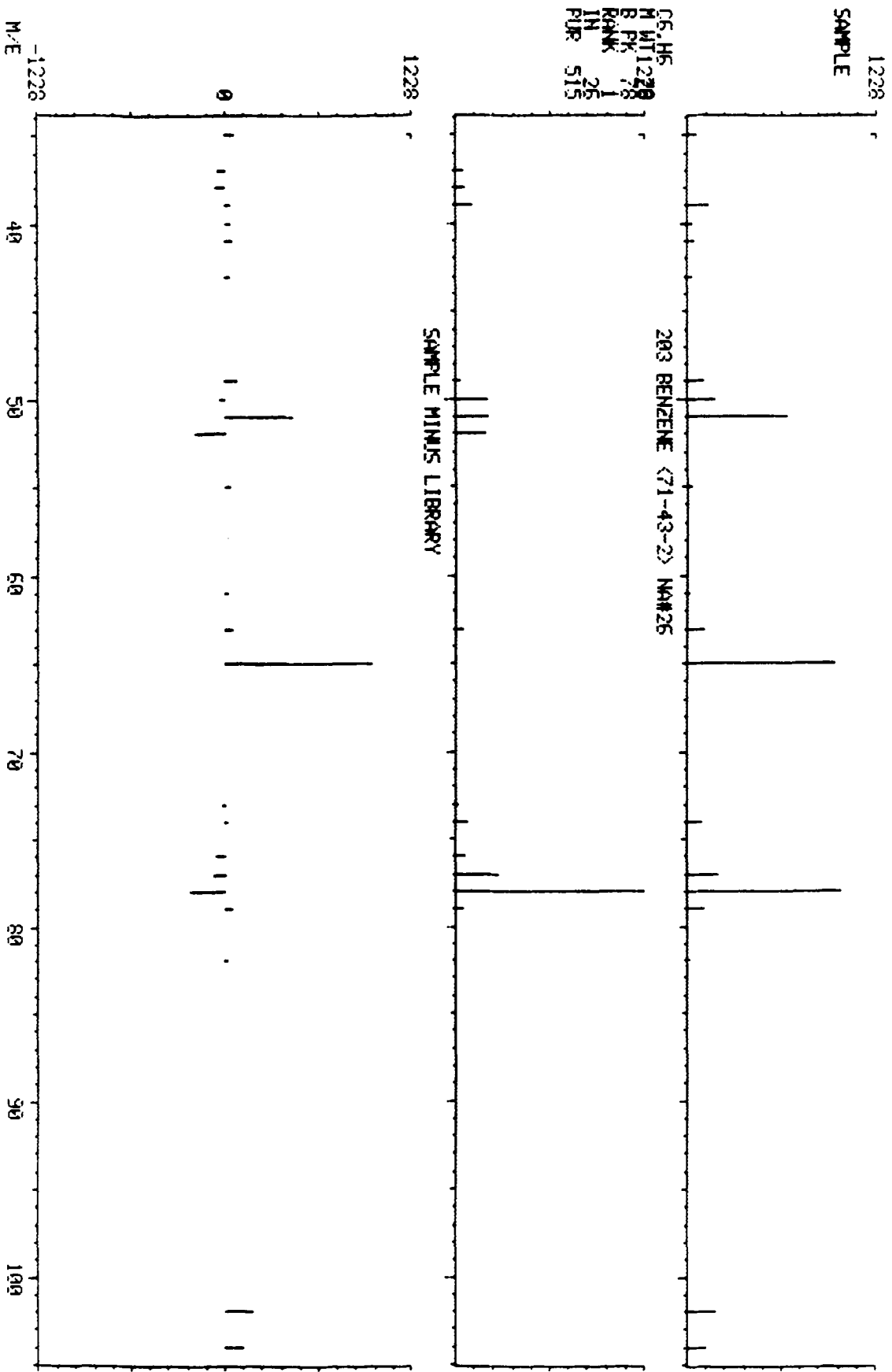
BASE M/E: 49/ 49
RIC: 8415.7 10607.



LIBRARY SEARCH
 12/29/89 1:52:00 + 5:48
 SAMPLE: 300L GC#309687 EPA#B202TAR RR CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

COMPUCHEM LABS
 DATA: C3R09687C13 # 464
 BASE M/E: 78
 RIC: 4223.

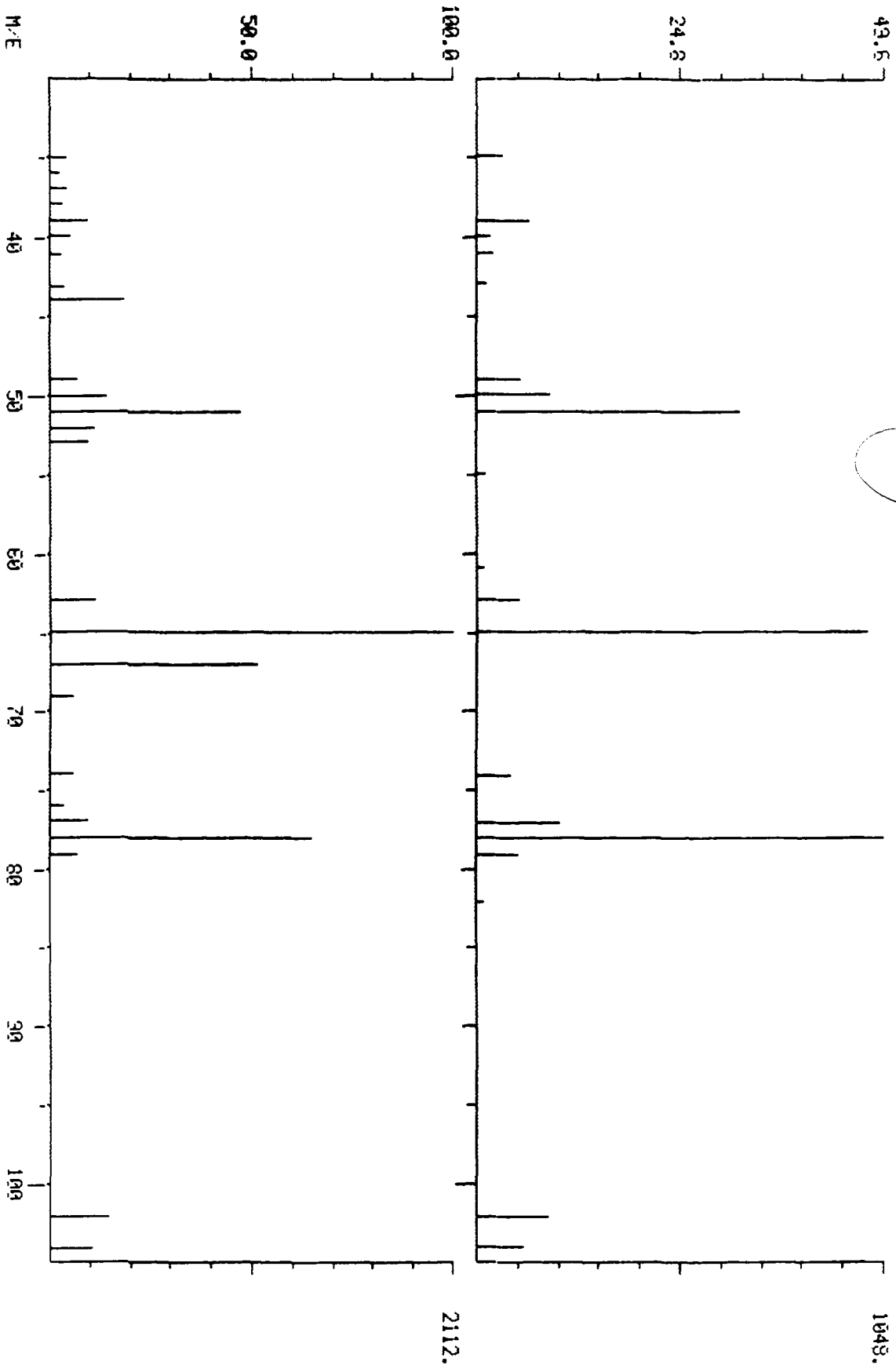
C6.H6
 MW 78
 B.P. 178
 BOILING 178
 IN 26
 PUR 515



DUAL MASS SPECTRUM
12/29/89 1:52:00 + 5:48
SAMPLE: 30UL CCM309587 EPA#B2021AR PE CASE#18756 ON #13
ENHANCED (5 158 2H) 293 BENZENE (71-43-2) HA#25

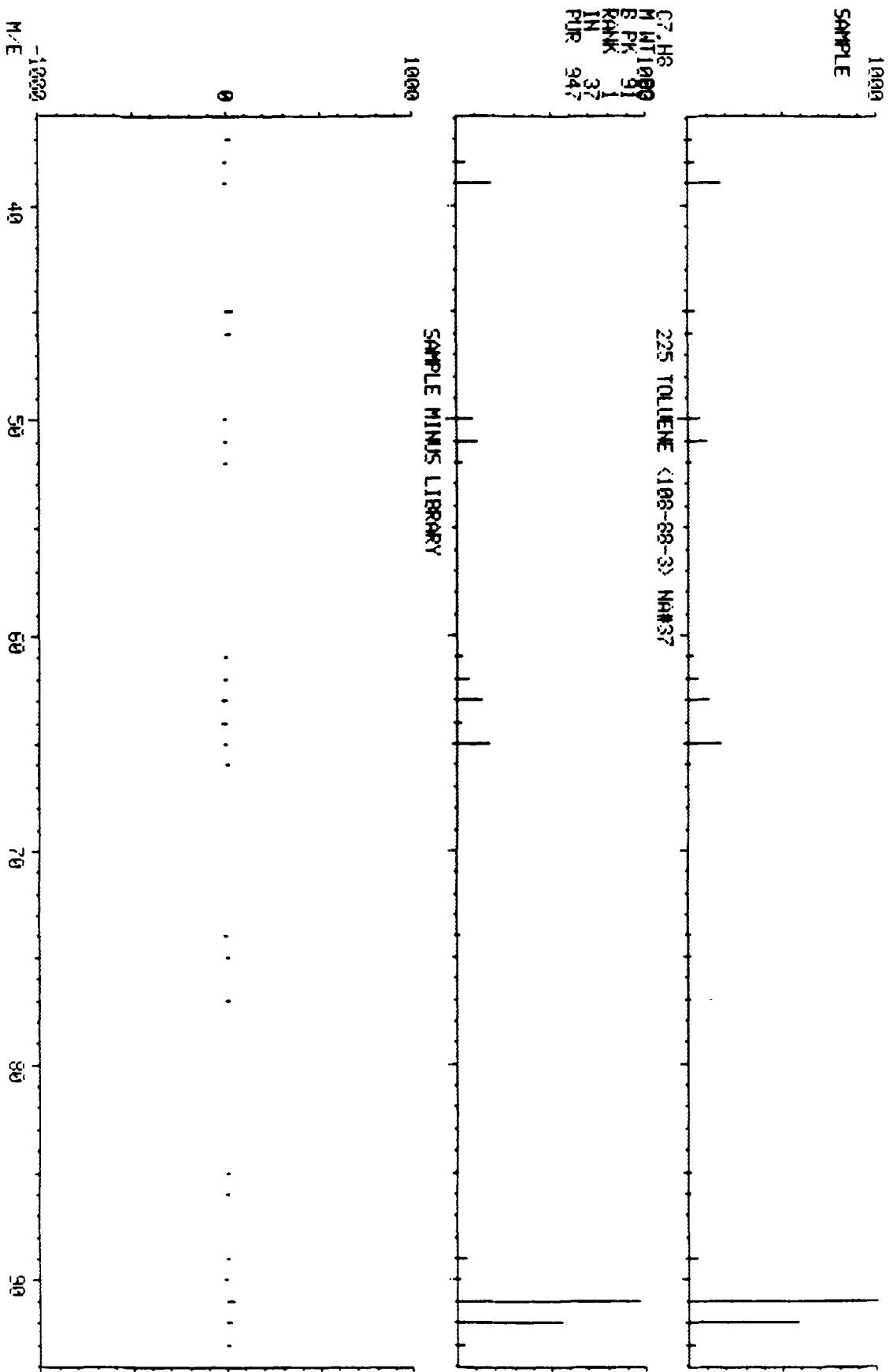
COMPUchem LABS
Dr. J. J. Van
1/1/90

DATA: C3R09587C13 #464
BASE M/E: 78/ 65
RIC: 4223.7/ 8975.



LIBRARY SEARCH
 12/29/89 1:52:00 + 8:54
 SAMPLE: 30UL CCM309687 EPA#B202TAR RE CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

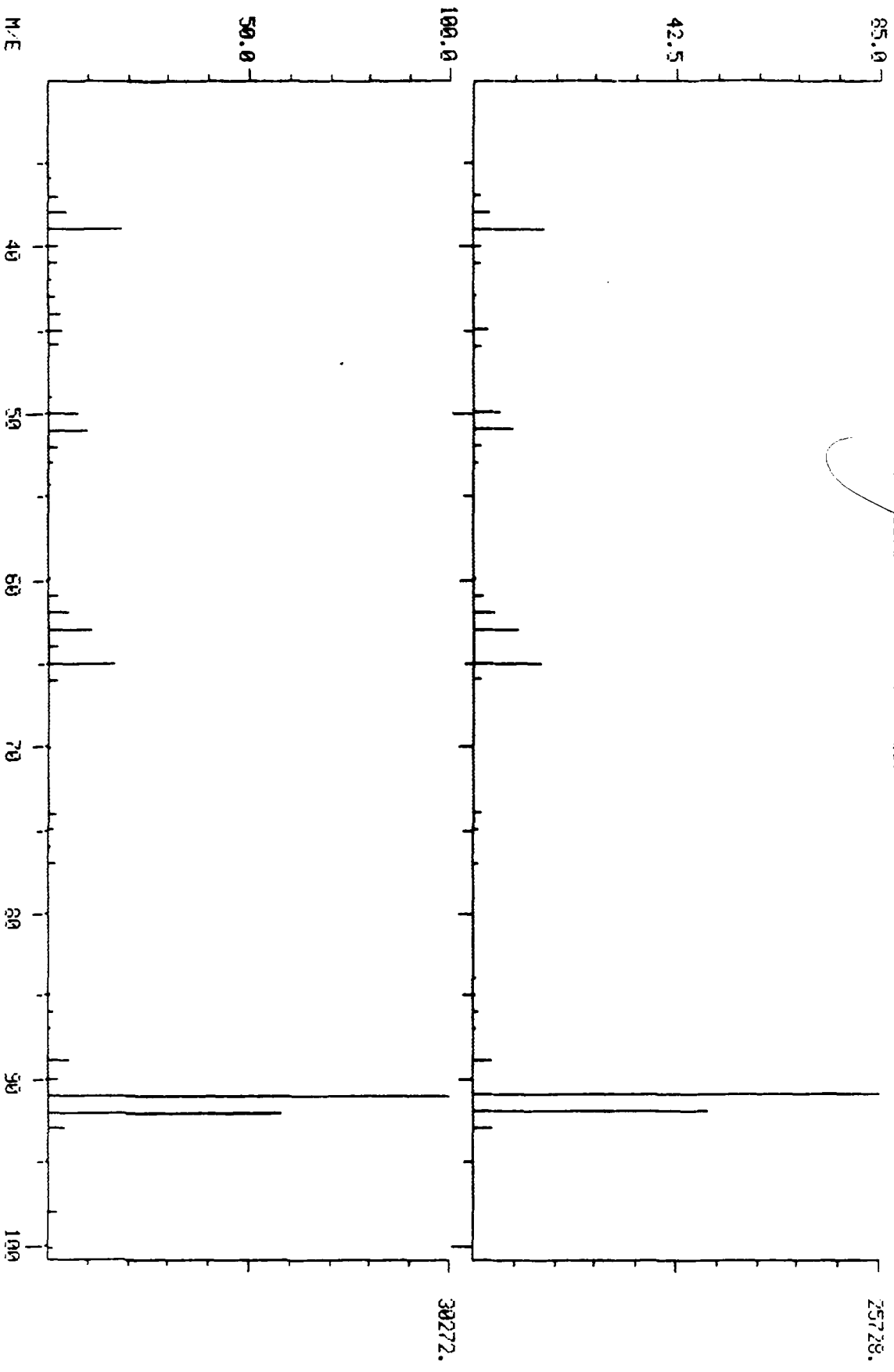
COMPUCHEM LABS
 DATA: CCR09687C13 # 712
 BASE M/E: 31
 RIC: 65535.



DUAL MASS SPECTRUM
12/29/89 1:52:00 + 8:54
SAMPLE: 30UL CC#309687 EPA#B2R21AR RE CASE#18756 ON #13
ENHANCED (S 158 2N)

COMPUchem LABS
DL 944 1340

DATA: C9F09687C13 #712 BASE M/E: 91/ 91
RIC: 67071. / 86271.

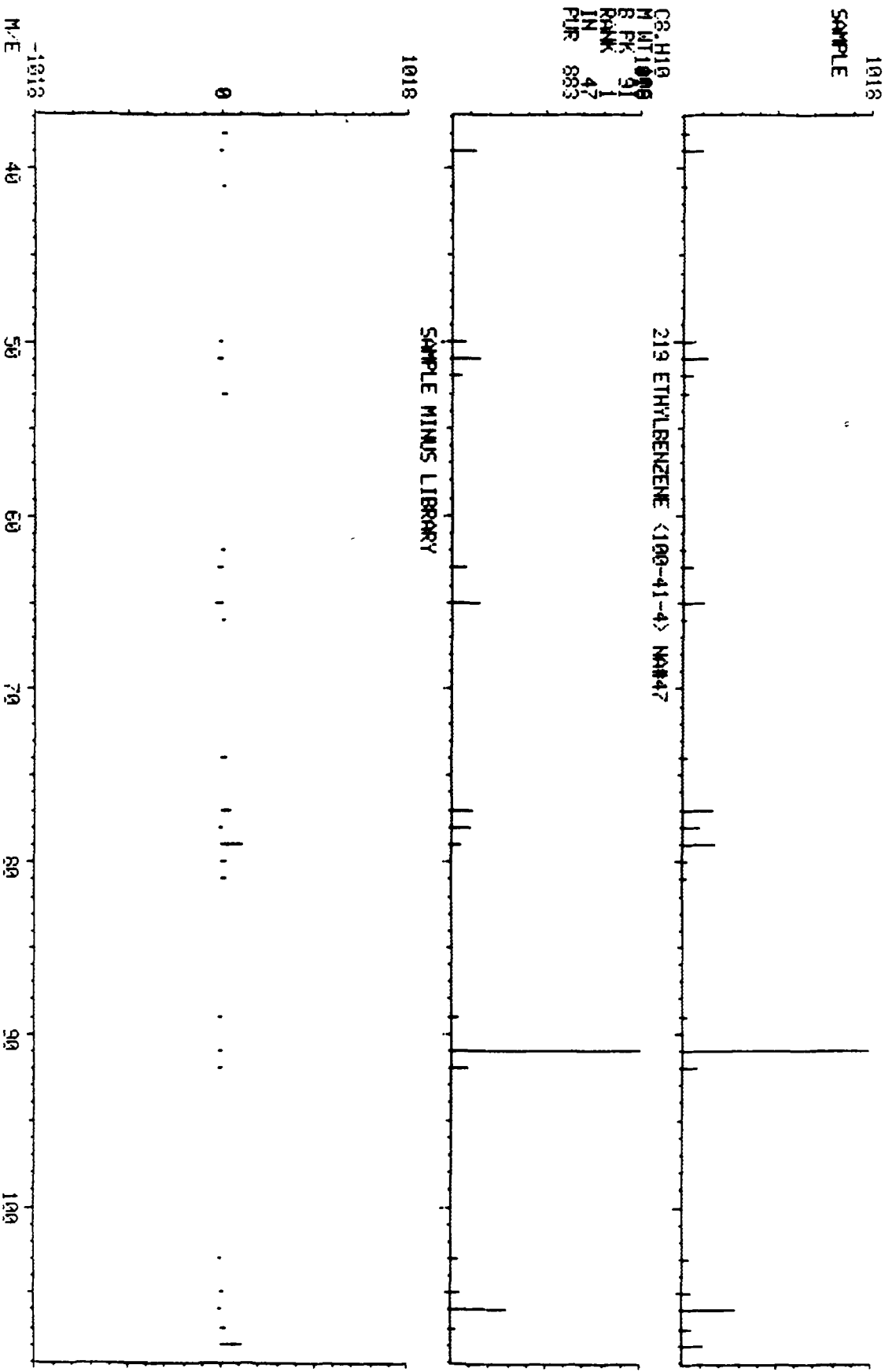


COMPUCHEM LABS
LIBRARY SEARCH
12/29/89 1:52:00 + 11:40
SAMPLE: 30UL CC#309687 EPA#B202TAR RE CASE#18756 ON #13
ENHANCED (5 158 2N 0T)

DL SAID 1/15/90
DATA: C309687C13 # 333

BASE M/E: 91
RIC: 182527.

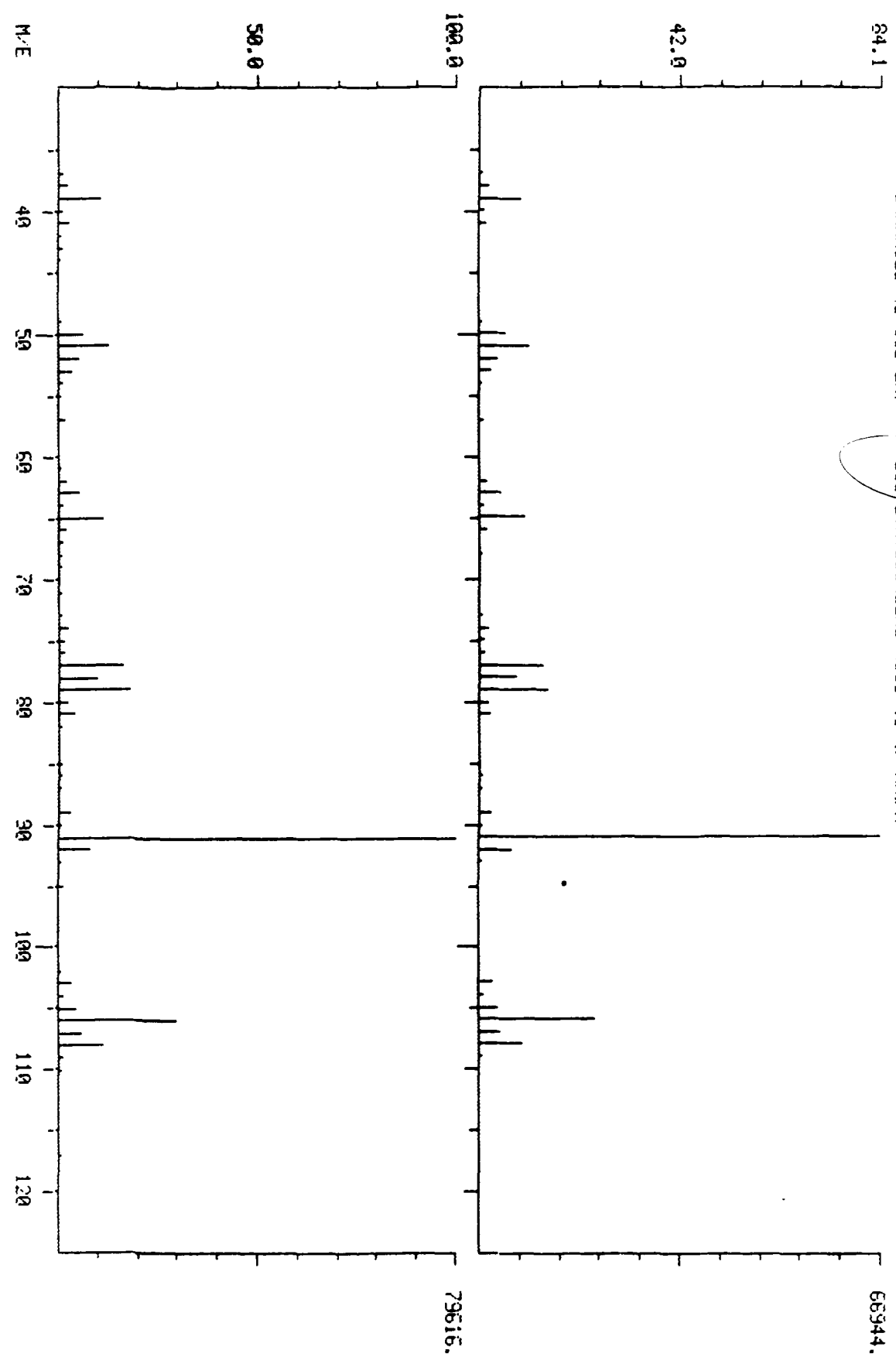
CG-H18
M UTI 100
B PK 91
IN 47
PUR 883



DUAL MASS SPECTRUM
12/23/89 1:52:00 + 11:40
SAMPLE: 30UL CC#309687
ENHANCED (S 158 2N)

CONFICHEM LABS
DATA: C3R09687013 #933
EPA#8202TAR PC CASE#18756 ON #13
219 ETHYLBENZENE <100-41-4> NA#47

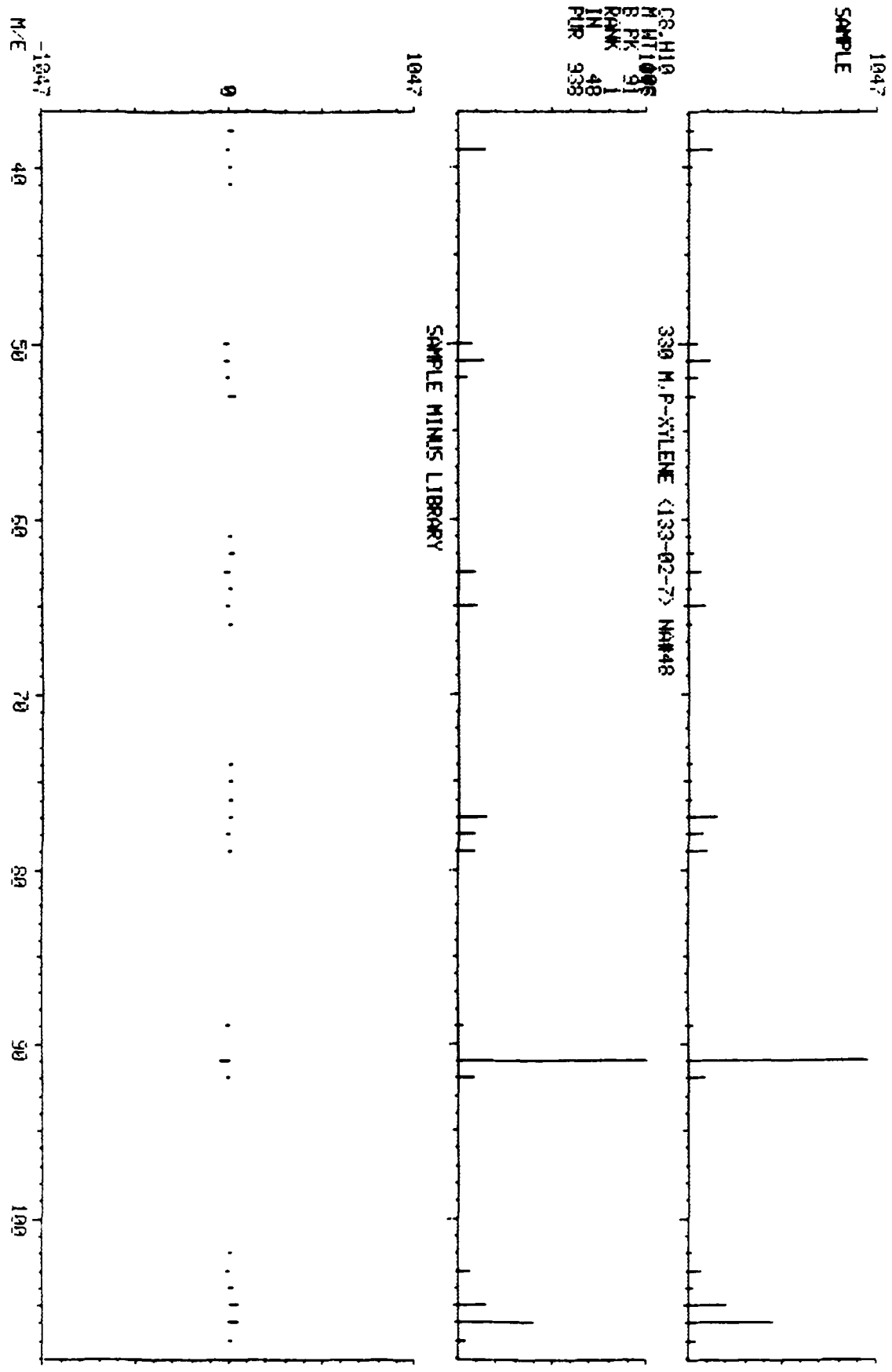
BASE M/E: 91 / 91
RIC: 191743. / 243967.



COMPUCHEN LABS
 12/29/89 1:52:00 + 11:54
 SAMPLE: 39UL CC#309687 EPA#B202TAR RZ CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

DATA: C3R09687013 # 952
 BASE M/E: 91
 RIC: 312831.

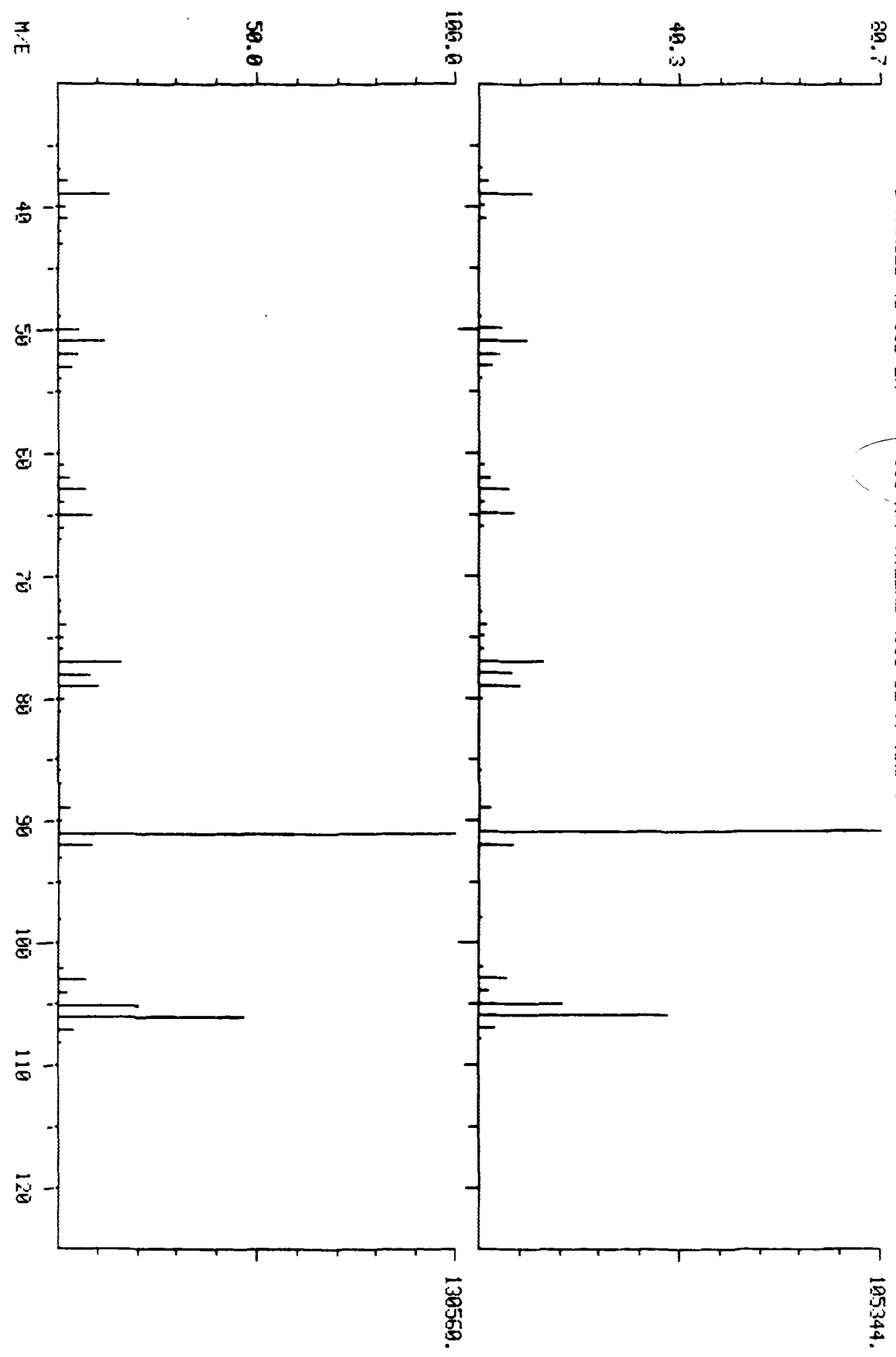
CG-H110
 M HT 1005
 R PK 91
 RANK 1
 IN 48
 PLR 938



DUAL MASS SPECTRUM
12/29/89 1:52:00 + 11:54
SAMPLE: 30UL C6H3095687
ENHANCED (S 158 2N)

COMPUCHER LABS
DATA: C3R095687C13 #952
RIC: 318463. / 401407.
330 M,P-XYLENE (133-02-7) NA#48

BASE M/E: 91 / 91
M/E: 318463. / 401407.

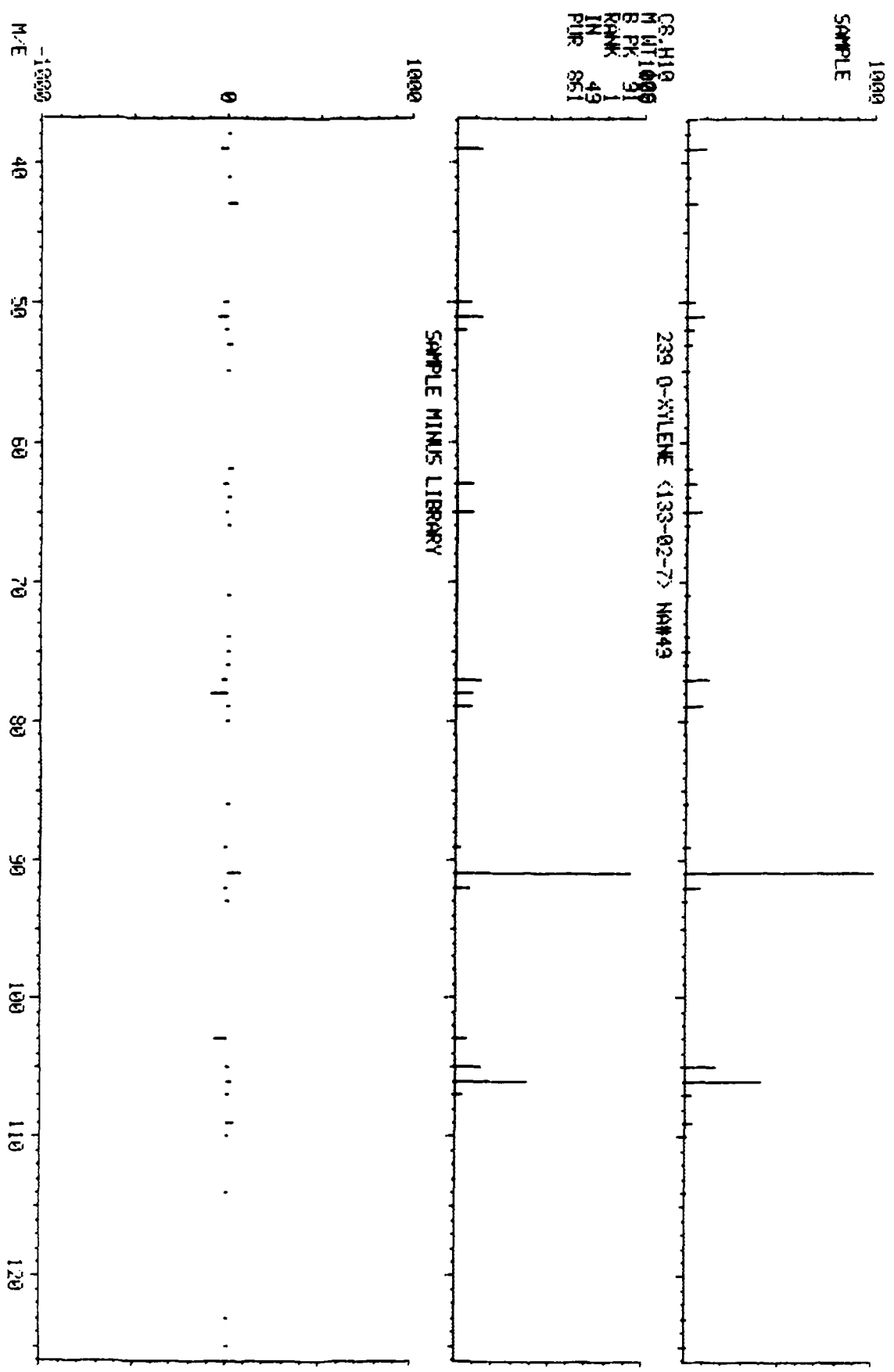


LIBRARY SEARCH
12/29/89 1:52:00 + 12:34
SAMPLE: 30UL CC#309687 EPA#B2021AR-RE CASE#18756 ON #13
ENHANCED (S 158 2N 0T)

COMPUchem LABS
DC 30UL/137D

DATA: C3R09687C13 #1006

BASE M/E: 31
PIC: 179967.

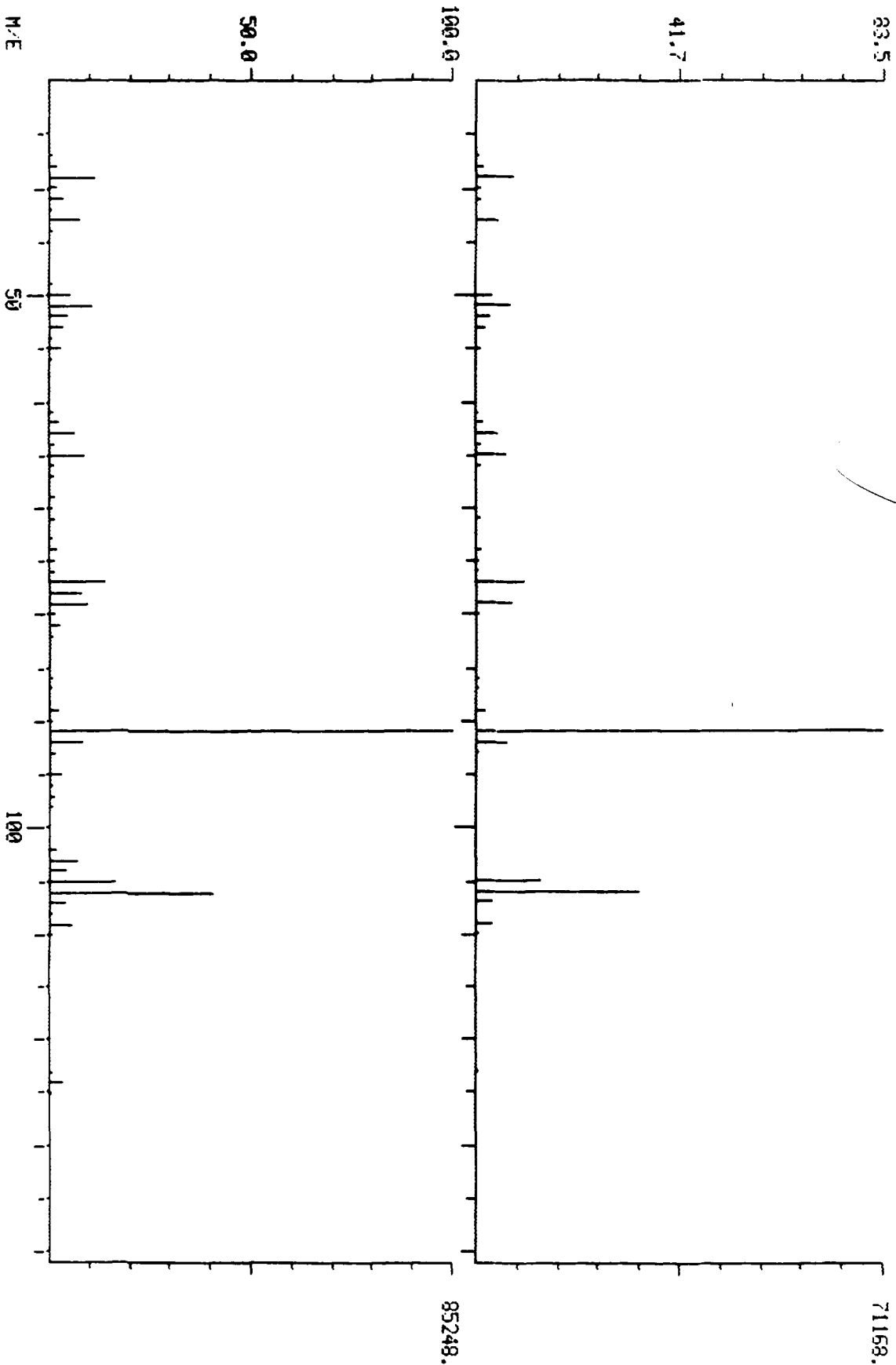


DUAL MASS SPECTRUM
12/29/89 1:52:00 + 12:34
SAMPLE: 30UL CCK309687 EPANB2021AR PE CASE#18756 ON #13
ENHANCED (5 158 2N)

DL 2008 1/14/90
COMPUchem LABS
239 D-XYLENE (133-02-7) NR#49

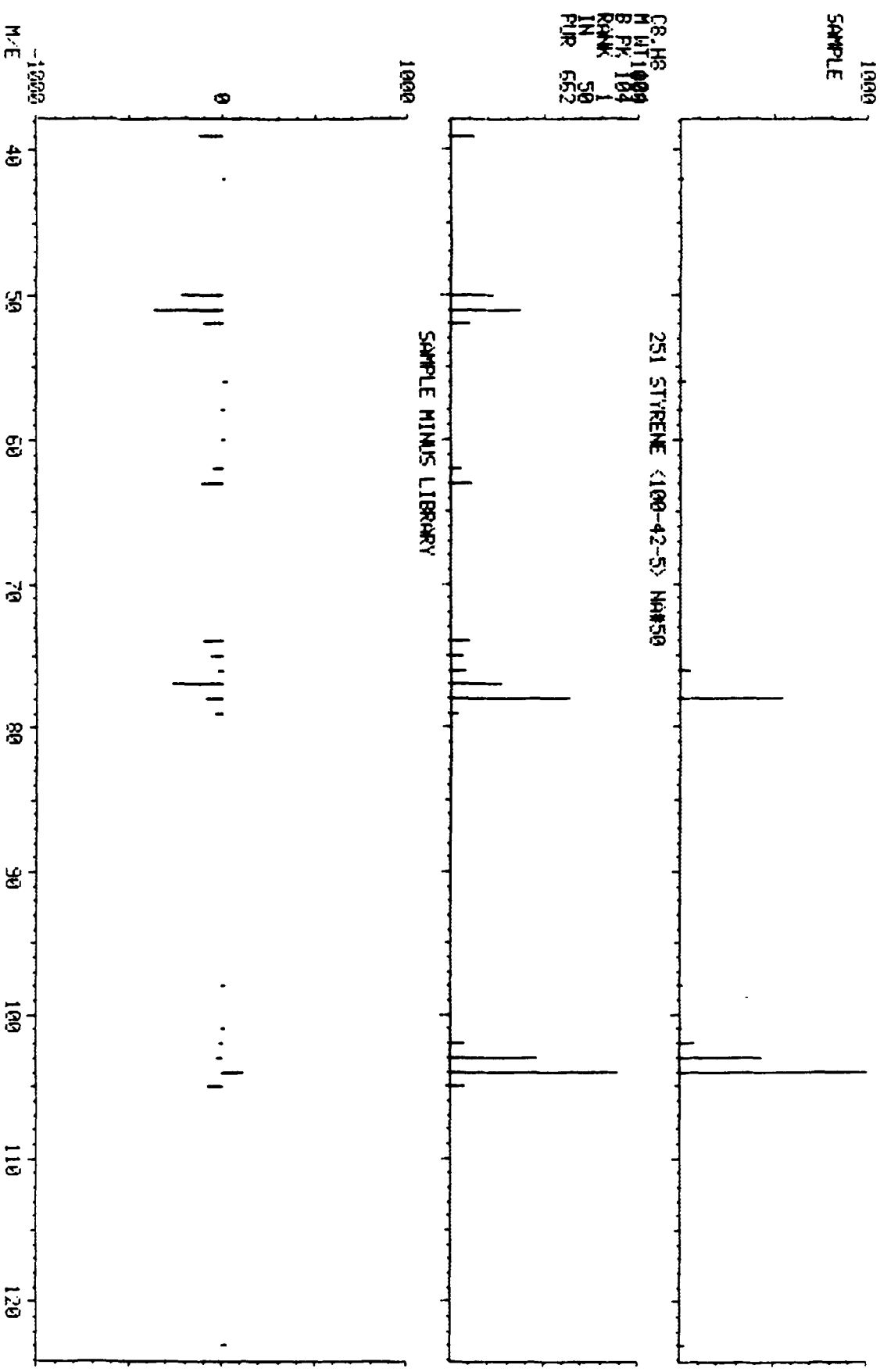
COMPUchem LABS

DATA: C3K09687013 #1006 BASE M/E: 91 / 91
RIC: 182527. / 274431.



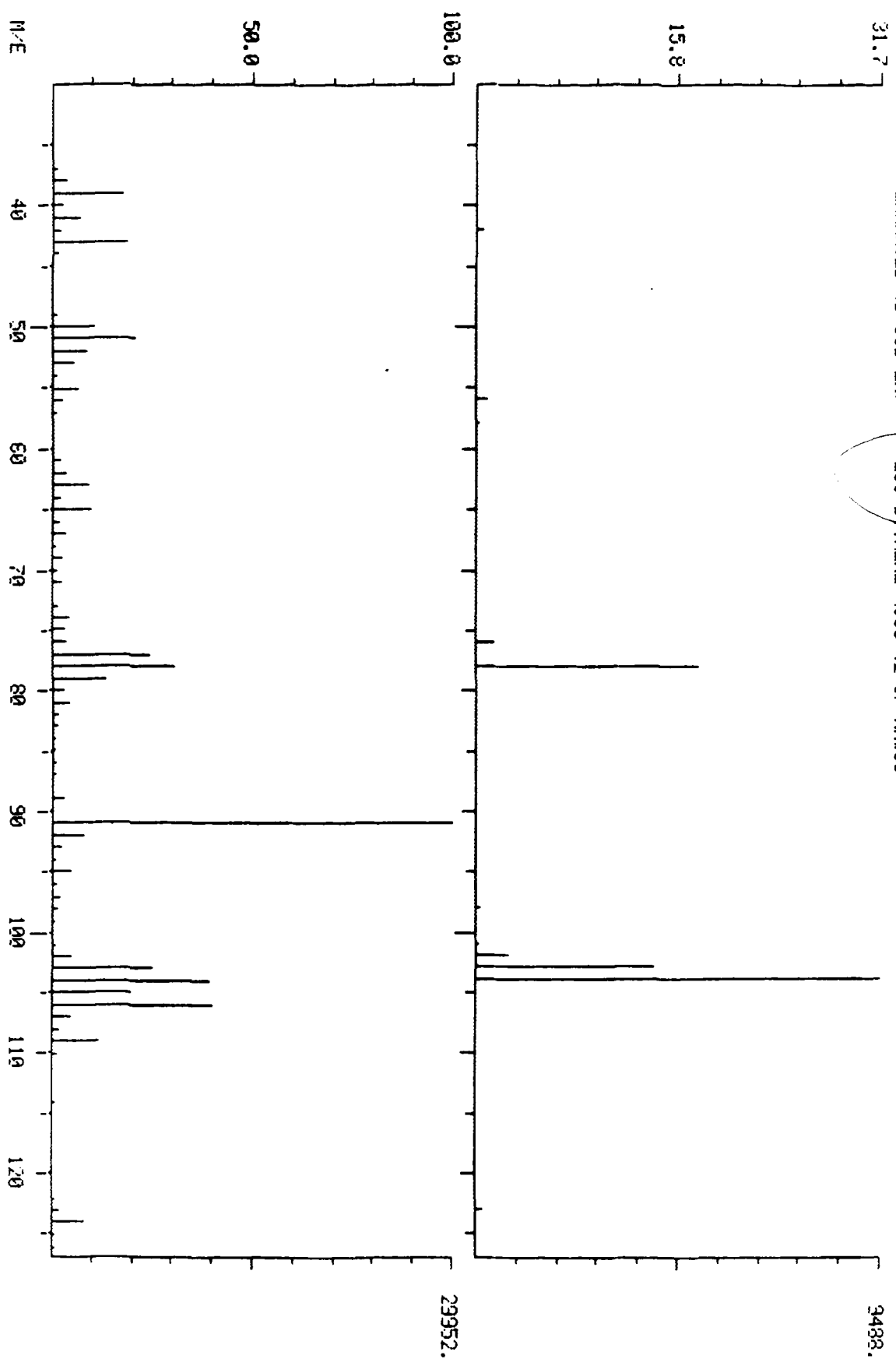
LIBRARY SEARCH
12/29/89 1:52:00 + 12:33
SAMPLE: 30UL OC#309687 EPA#82021AR RE CASE#18756 ON #13
ENHANCED (5 158 2H QT)

COMPUchem LABS
DL SA# 113/90
DATA: CAR09687C13 #1012
BASE M/E: 104
RIC: 20863.



DUAL MASS SPECTRUM
12/29/89 1:52:00 + 12:39
SAMPLE: 380L CC#369687 EPA#B2A2TAR PE CASE#18756 ON #13
ENHANCED (S 158 2N) 251 STYRENE (100-42-5) NA#50

COMPUCHEN LABS
DATA: C3R09687C13 #1012 BASE M/E: 104 / 91
RIC: Z0863. / 154111.
3488.
29952.

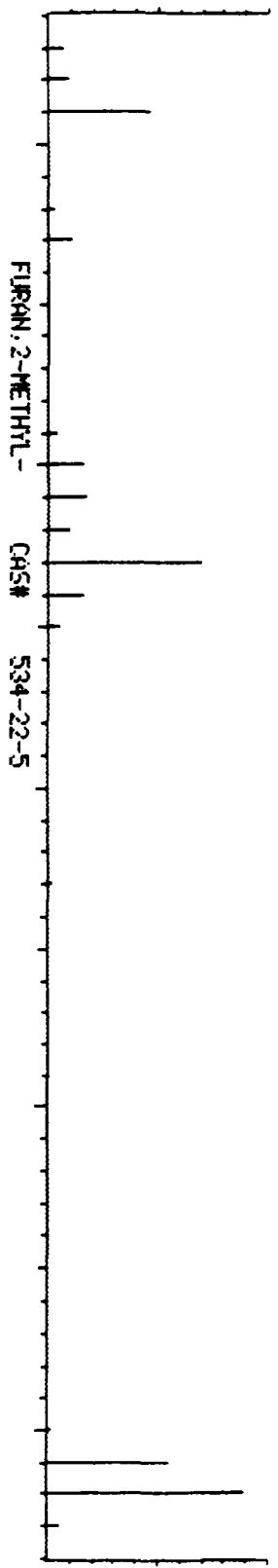


LIBRARY SEARCH
 12/29/89 1:52:00 + 4:13
 SAMPLE: 30UL CC#309687 EPA#B202TAR PE CASE#18756 ON #13
 ENHANCED (5 158 2N 0T)

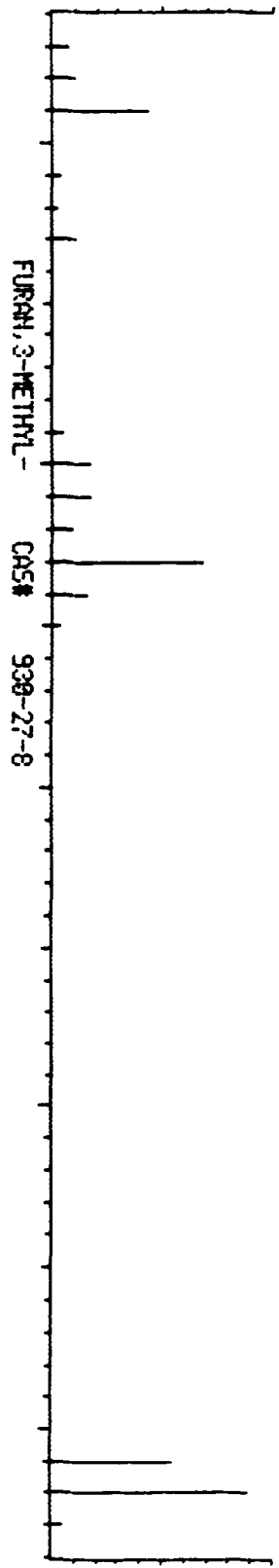
COMPUCHEM LABS
 DATA: C3R09687C13 # 337
 BASE M/E: 82
 RIC: 20383.

1134
SAMPLE

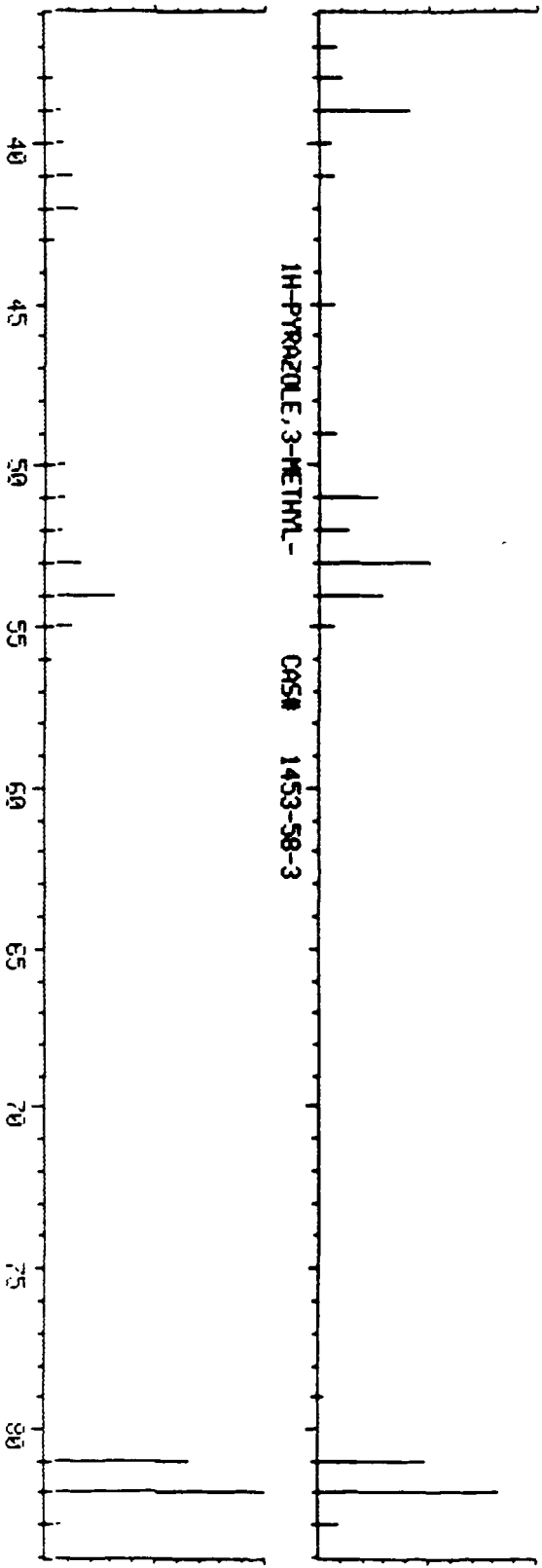
C5.H6.0
 M WT 1134
 B PK 82
 RANK 1
 IN 341
 PUR 965



C5.H6.0
 M WT 1134
 B PK 82
 RANK 342
 IN 876
 PUR 876



C4.H6.N2
 M WT 1134
 B PK 82
 RANK 3
 IN 340
 PUR 824

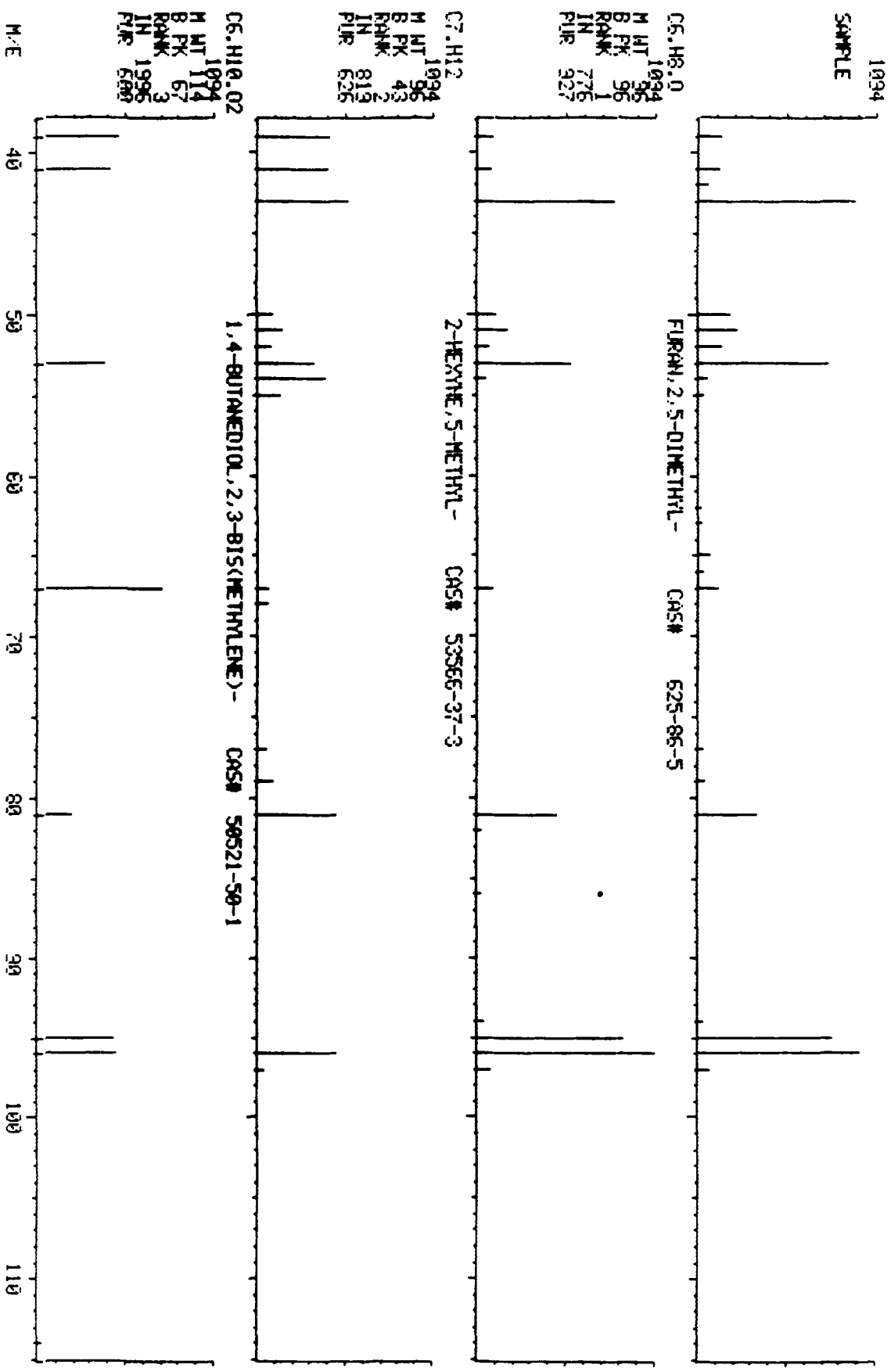


M/E

40 45 50 55 60 65 70 75 80

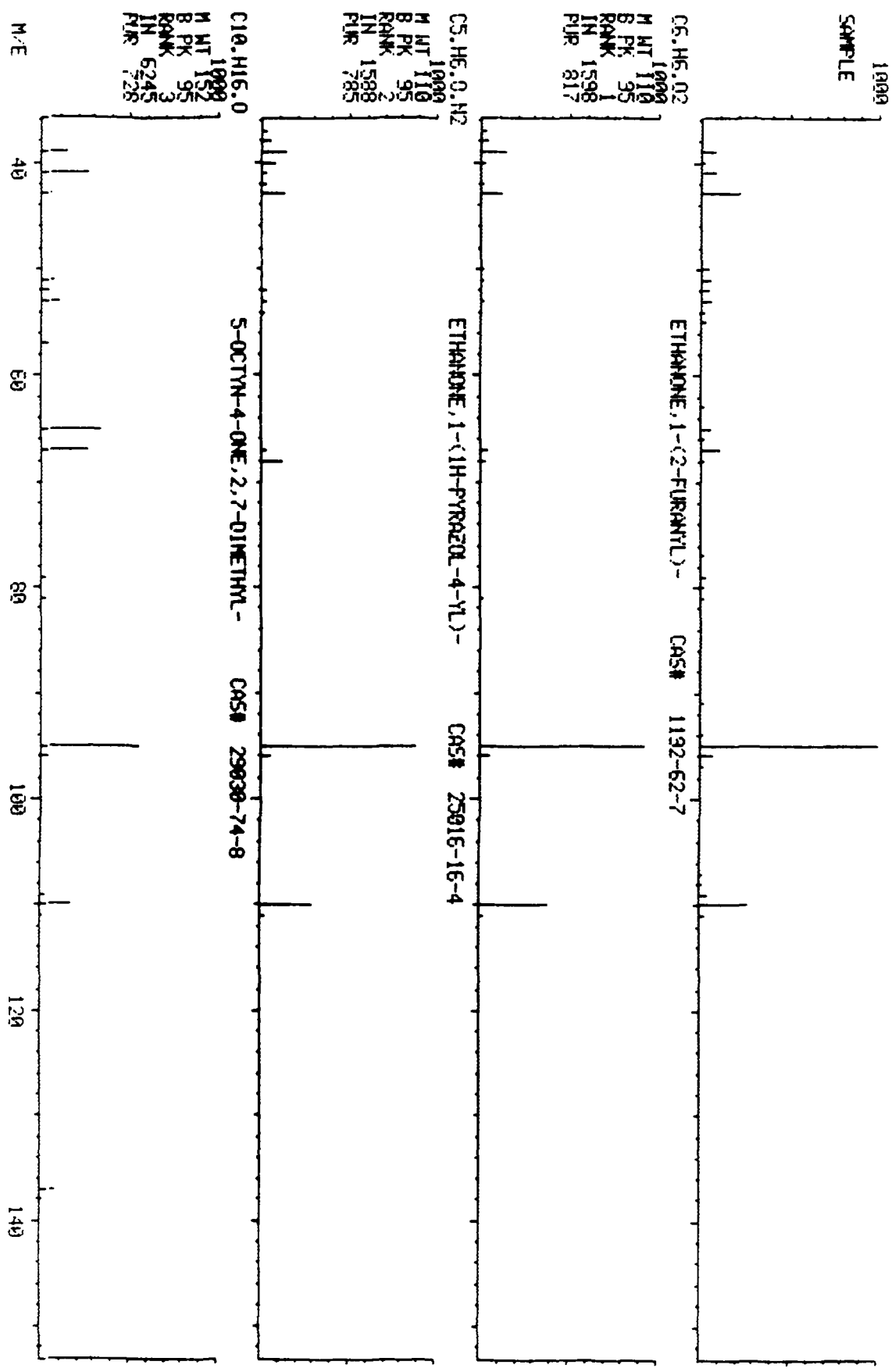
LIBRARY SEARCH
 12/29/89 1:52:00 + 7:10
 SAMPLE: 30UL CC#389687 EPA#B202TAR RE CASE#18756 ON #13
 ENHANCED (S 158 2H QT)

COMPUchem LABS
 DATA: C3R09687013 # 573
 BASE M/E: 95
 RIC: 39167



LIBRARY SEARCH
 12/23/83 1:52:00 + 9:50
 SAMPLE: 30UL CC#389687 EPA#8202TAR-RE CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

COMPUCHEN LABS
 DATA: C3R09687013 # 787
 BASE M/E: 95
 PIC: 84607.



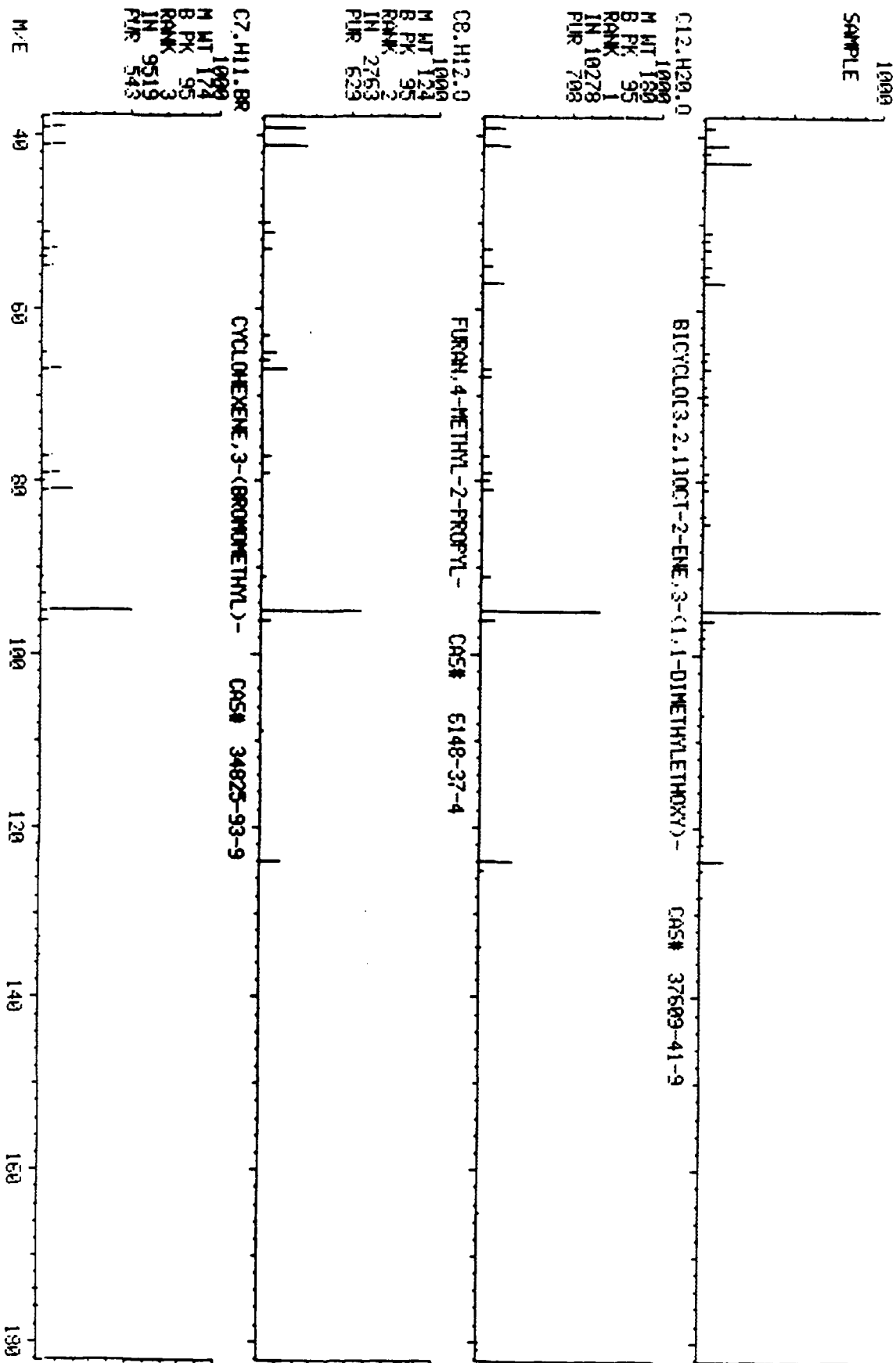
LIBRARY SEARCH
12/29/89 1:52:00 + 12:14
SAMPLE: 30UL CC#309587 EPA#B202TAR RE CASE#18756 ON #13
ENHANCED (S 158 2H 0T)

COMPUchem LABS

DL said 1/2/90

DATA: C3R09587C13 # 979

BASE M/E: 95
PIC: 99967.



LIBRARY SEARCH
 12/29/89 1:52:00 + 12:49
 SAMPLE: 39UL CC#309687 EPA#B2021AR PE CASE#18756 ON #13
 ENHANCED (5 158 21 0T)

COMPUCHEM LABS
 DATA: C3R09687C13 #1025
 BASE M/E: 109
 RIC: 55935.

1000
 SAMPLE

08.H12.0
 1000
 M UT 124
 B PK 109
 RANK 2771
 IN 761
 PUR 761

2-CYCLOPENTEN-1-ONE,3,5,5-TRIMETHYL- CAS# 24156-95-4

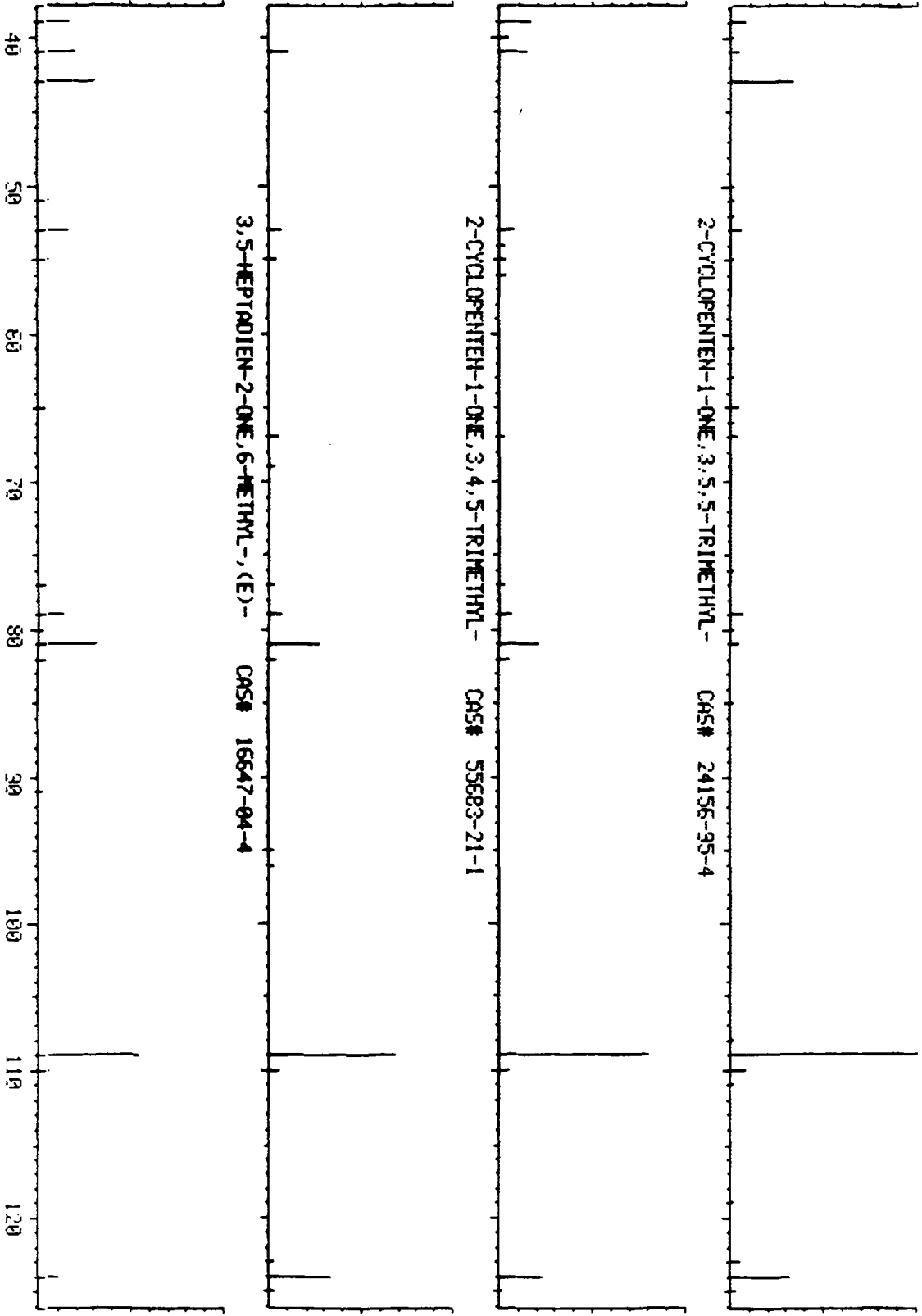
08.H12.0
 1000
 M UT 124
 B PK 109
 RANK 2780
 IN 754
 PUR 754

2-CYCLOPENTEN-1-ONE,3,4,5-TRIMETHYL- CAS# 55683-21-1

08.H12.0
 1000
 M UT 124
 B PK 109
 RANK 2769
 IN 716
 PUR 716

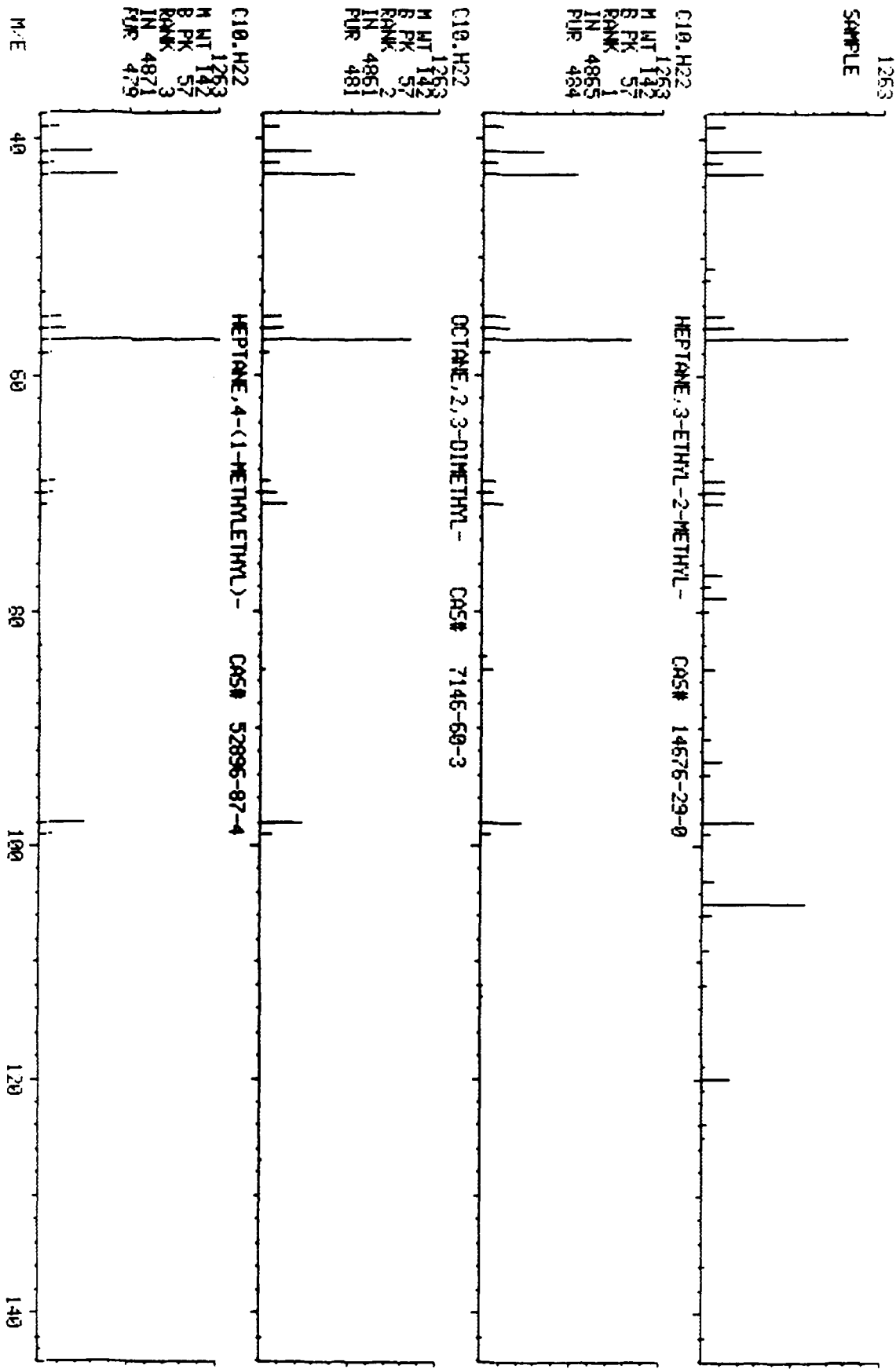
3,5-HEPTADIEN-2-ONE,6-METHYL-,(E)- CAS# 16647-04-4

M/E



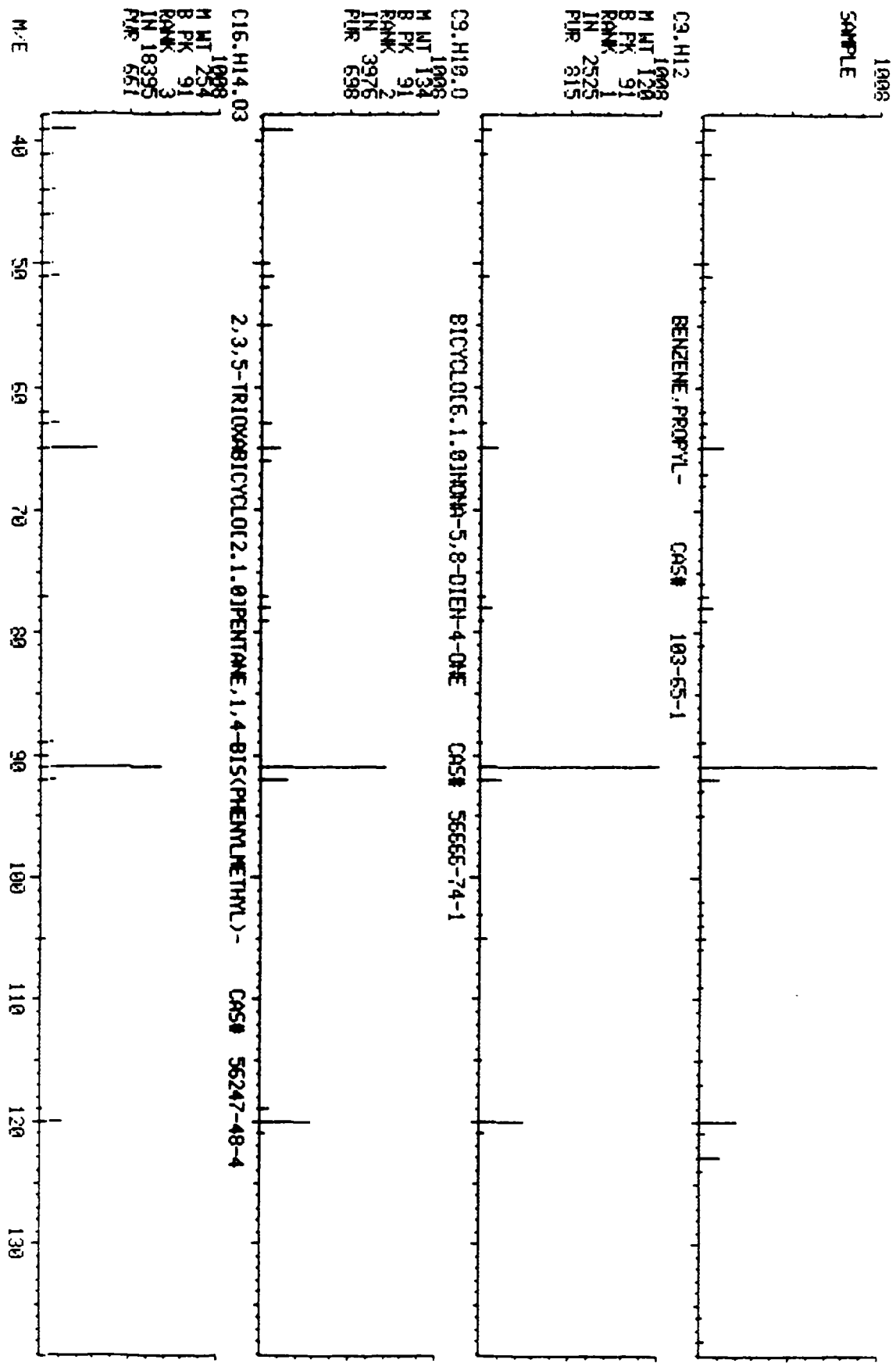
LIBRARY SEARCH
12/29/89 1:52:00 + 13:19
SAMPLE: 30UL CC#309587 EPANR202TAR-RE CASE#18756 ON #13
ENHANCED (S 158 2N 0T)

COMPUchem LABS
DATA: C3R09587013 #1065
BASE M/E: 57
RIC: 37695



LIBRARY SEARCH
 12/29/89 1:52:00 + 14:07
 SAMPLE: 30UL CC#309687 EPA#B202TAR RE CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

COMPUchem LABS
 DATA: C3P09687C13 #1129
 BASE M/E: 91
 RIC: 32415



LIBRARY SEARCH
12/29/89 1:52:00 + 14:21
SAMPLE: 30UL CC#309687 EPA#B202TAR RE CASE#18756 ON #13
ENHANCED (5 158 2N 0T)

COMPUCHEN L985
DL 204, 1570

DATA: C3R09687C13 #1148

BASE M/E: 105
RIC: 222463.

1124
SAMPLE

C9.H12
M WT 1124
B PK 105
RANK 1
IN 2528
PIR 856

BENZENE, 1-ETHYL-2-METHYL- CAS# 611-14-3

C9.H12
M WT 1124
B PK 105
RANK 2
IN 2529
PIR 845

BENZENE, 1-ETHYL-3-METHYL- CAS# 620-14-4

C9.H12
M WT 1124
B PK 105
RANK 3
IN 2524
PIR 844

BENZENE, (1-METHYLETHYL)- CAS# 98-82-8

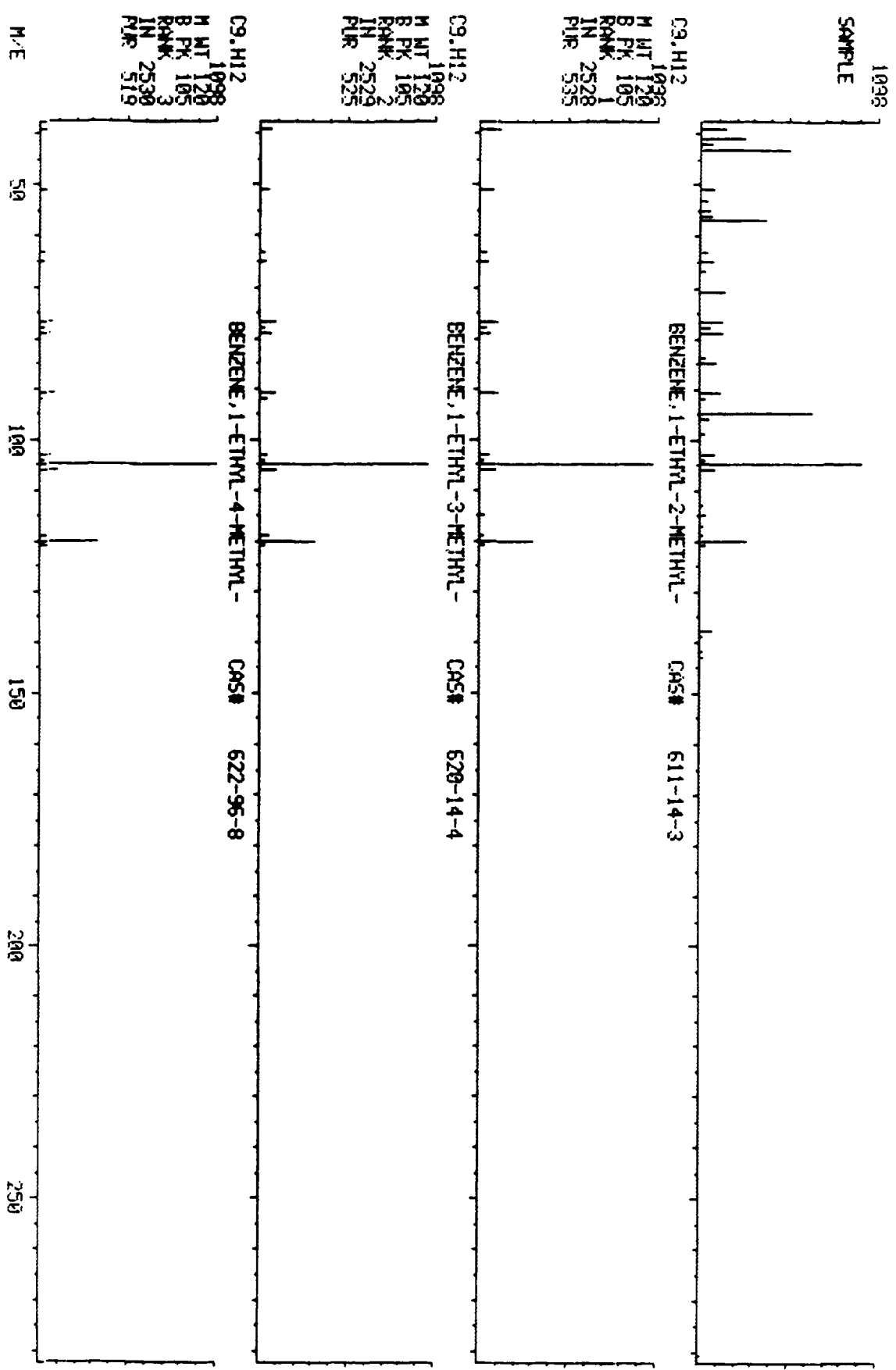
M/E 40 50 80 100 120 140 160

LIBRARY SEARCH
 12/29/83 1:52:00 + 14:49
 SAMPLE: 30UL CC#309587 EPA#B202TAR RE CASE#18756 ON #13
 ENHANCED (S 158 2N 0T)

COMPUchem LABS

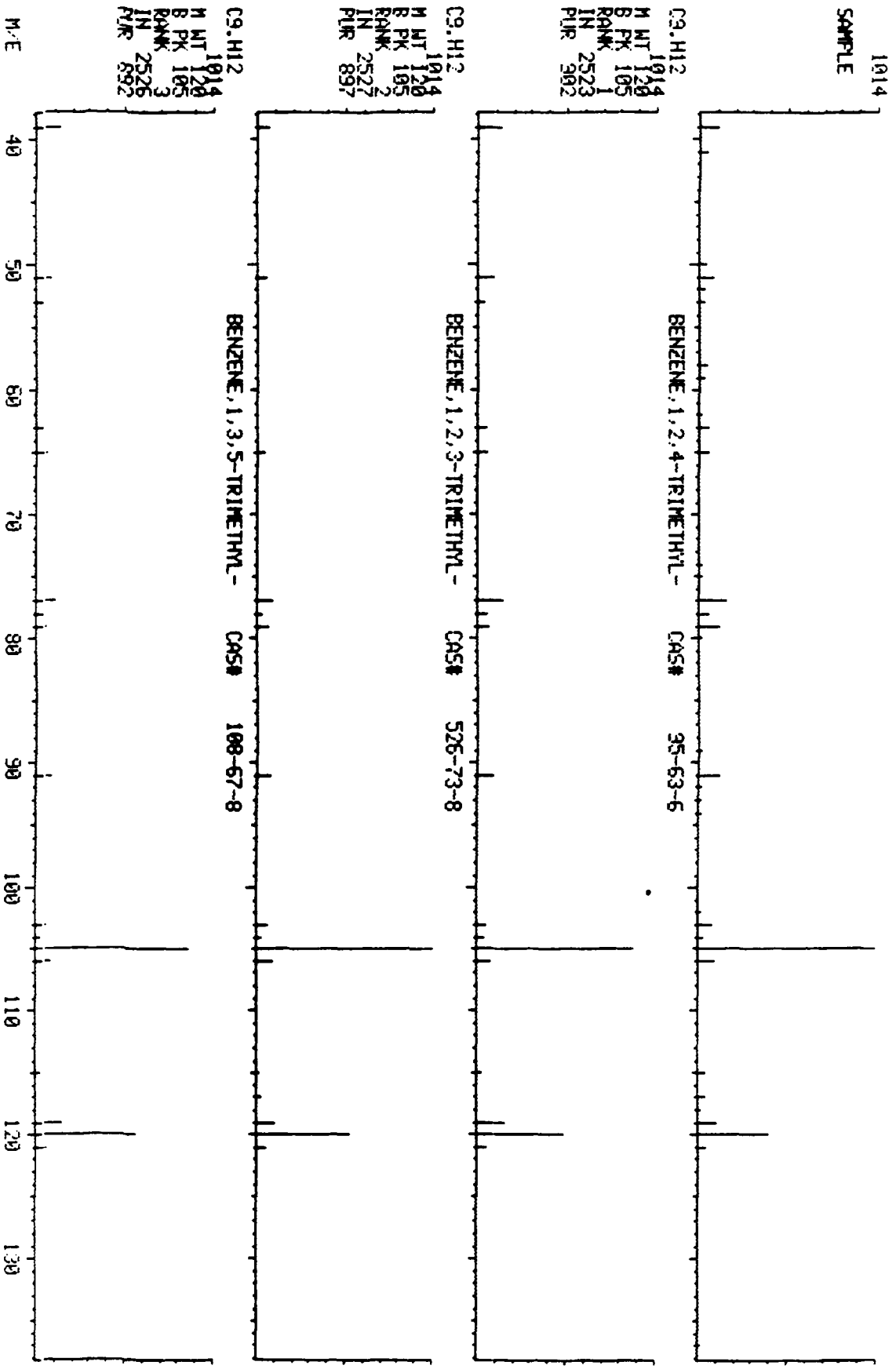
DATA: C3R0687C13 #1186

BASE M/E: 105
 RIC: 139335



LIBRARY SEARCH
 12/29/89 1:52:00 + 15:11
 SAMPLE: 30UL CC#309687 EPA#8202TAR RE CASE#18756 ON #13
 ENHANCED (5 158 2N 0T)

COMPUchem LABS
 DL *sa* # 157D DATA: C3R09687C13 #1215
 BASE M/E: 105
 RIC: 259923.



AB INSTRUCTIONS:

RECEIPT DATE

CASE#: 18756

DUE DATE

VOA
C/MS WORKSHEET

COMPUCHENO: 309687R³

RE R2E D (:)
R3E R4E D2E (:)

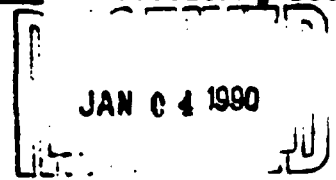
MED. SOLID, EPA 80W 2/87

Sample Prep Code---156
Instrument Code---414
Compound List---495
Surrogate Std-----394
Internal Std-----036

SAMPLE ID: 1370ZTARRE Dry Wt. Factor 2.78 X Moisture 64

GC/MS ANALYSIS

Amount Purged: 10mls soil or Dilution 30 ul/10000ul/Xg soil
Internal Standard Volume Added 5 ul
Surrogate Standard Volume Added 5 ul
FB Filename 136891228A14 Disk ()
Blank Filename CAD11246B13 Disk ()
Standard Filename CTR91228A13 Disk ()
Sample Filename C3R09687R13 Disk ()



ANALYST(S): Injection 1452 N. H. H. Work-up 11-2 N. H. H.

GC/MS REVIEW

CONDITION
CODE

DA

Entry Codes OK, EA, ES, SH, JS, SL, SH, JA, DA

Non-Entry Codes IM, IL, IH, SU, CT, CS, PC, NR
IF, LA, DI, CO, RN, DU, SI, SF
UP, BB, OT, VC, FO, NS



Disposition: Complete
 Reprep neat required
 Reprep using _____
 Dilute (:)

Extraneous Peak Search Results:

of Peaks Found: 10

Quality Assurance Notice(s):

Notices Required 1

COMMENTS:

GC/MS Review [Signature] Date 1/3/90 Auditor _____ Date ____/____/____

REPORT INTEGRATION

Final Reportable Package(s): C3R09687B13 / G2R09687C17 Total # of Injections: _____

QA COMMENTS:

Initials _____ Date ____/____/____

FINAL REVIEW:

Initials _____ Date ____/____/____

AC1010 (4/89)

CMP #	M/E F	COMPOUND NAME	SCAN	AREA	QUANT REPORT VALUE	REPORTED AMOUNT (UG/KG)	DETECT LIMIT (UG/KG)
234	128 I	BROMOCHLOROMETHANE (IS)	397	49600	50.0		2100
221	50	CHLOROMETHANE				BDL	4200
231	62	VINYL CHLORIDE				BDL	4200
220	94	BROMOMETHANE				BDL	4200
209	64	CHLOROETHANE				BDL	4200
216	96	1,1-DICHLOROETHENE				BDL	2100
254	76	CARBON DISULFIDE				BDL	2100
252	43	ACETONE (2-PROPANONE)			13.2	BDL	4200
248	114 I	1,4-DIFLUOROBENZENE (IS)	530	173000	50.0		2100
222	84	METHYLENE CHLORIDE			14.1	15000	2100
226	96	TRANS-1,2-DICHLOROETHENE				BDL	2100
214	63	1,1-DICHLOROETHANE				BDL	2100
257	43	VINYL ACETATE				BDL	4200
237	96	CIS-1,2-DICHLOROETHENE				BDL	2100
253	72	2-BUTANONE				BDL	4200
211	83	CHLOROFORM				BDL	2100
227	97	1,1,1-TRICHLOROETHANE				BDL	2100
206	117	CARBON TETRACHLORIDE				BDL	2100
203	78	BENZENE			4.3	5000	2100
215	62	1,2-DICHLOROETHANE				BDL	2100
270	117 I	D5-CHLOROENZENE (IS)	906	250000	50.0		2100
229	130	TRICHLOROETHENE				BDL	2100
217	63	1,2-DICHLOROPROPANE				BDL	2100
212	83	BROMODICHLOROMETHANE				BDL	2100
218	75	CIS-1,3-DICHLOROPROPENE				BDL	2100
256	43	4-METHYL-2-PENTANONE			2.4	10000	4200
225	92	TOLUENE			91.3	110000	2100
250	75	TRANS-1,3-DICHLOROPROPENE				BDL	2100
228	97	1,1,2-TRICHLOROETHANE			1.9	79000	2100
224	164	TETRACHLOROETHENE				BDL	2100
255	43	2-HEXANONE				BDL	4200
208	129	DIBROMOCHLOROMETHANE				BDL	2100
207	112	CHLOROBENZENE				BDL	2100
219	106	ETHYLBENZENE			90.4	160000	2100
330	106	M, P-XYLENE			148.0	170000	2100
239	106	O-XYLENE			98.9	110000	2100
251	104	STYRENE			19.5	23000	2100
205	173	BROMOFORM				BDL	2100
223	83	1,1,2,2-TETRACHLOROETHANE				BDL	2100
258	65 S	D4-1,2-DICHLOROETHANE NA#57			16.5	110.0%	
247	95 S	BROMOFLUOROBENZENE			16.6	111.0%	
233	98 S	D8-TOLUENE NA#59			12.4	83.0%	
289	106	XYLENES (TOTAL)			247.0	100000	2100

CORRECTED/REVIEWED BY

DATE

(GC/MS DATA REVIEWER)

VOLATILE - MEDIUM LEVEL SOLID

CMP					QUANT	REPORTED	DETECT.	
#	M/E	F	COMPOUND NAME	SCAN	AREA	VALUE	AMOUNT	LIMIT
							(UG/KG)	(UG/KG)
299	96		1,2-DICHLOROETHENE (TOTAL)				BDL	2100
CHECKSUMS:								
			3979.	1833	472600.	913.3	296894.	

CORRECTED/REVIEWED BY


(GC/MS DATA REVIEWER)

DATE

12/31/67

NO	CC ID#	SURROGATE COMPOUND	QUANT REPORT VALUE	QUANT REPORT AMOUNT SPIKED	% ++ RECOVERY	CONTROL RANGE	P	F
40	258	D4-1,2-DICHLOROETHANE NA#57	16.5	15.0	110.	70-121	X	
41	247	BROMOFLUOROBENZENE	16.6	15.0	111.	74-121	X	
42	233	D8-TOLUENE NA#59	12.4	15.0	83.	81-117	X	

* ADVISORY SURROGATE ONLY

++ % RECOVERY = QUANT REPORT VALUE / QUANT REPORT AMOUNT SPIKED X 100 %

CORRECTION FACTOR CALCULATION:

$$\frac{4.0 \text{ G}}{\text{WET WEIGHT OF SAMPLE (G)}} \times \text{LIBRARY ADJUSTMENT} \times \text{DRY WEIGHT FACTOR} \times \frac{10000 \text{ UL}}{\text{UL USED}} =$$

$$\frac{4.0 \text{ G}}{4.00 \text{ (G)}} \times 1.25 \times \frac{2.78}{1.00} \times \frac{10000 \text{ UL}}{30. \text{ UL}} = 417.000$$

QUANT REPORT AMOUNT SPIKED CONVERSION FACTOR:

$$\frac{100. \text{ UL}}{\text{UL USED}} =$$

$$\frac{100. \text{ UL}}{30. \text{ UL}} = 3.3$$

VERSION 8

CORRECTED/REVIEWED BY

(GC/MS DATA REVIEWER)

DATE

[Handwritten Signature]
 12/09/07

LABORATORY NOTICE
COMPUCHEM ID # 309687
CLIENT ID # 8202TAR
CASE # 18756

The volatile fraction of this sample was prepared and analyzed as both a Low Level Solid and a Medium Level Solid. The initial analysis of 1 g sample from the Low Level preparation indicated that a Medium Level analysis was necessary in order to bring target compound concentrations into the instrumental analytical range. The analysis of 30 ul of the Medium Level extract proved to be an acceptable dilution. However, the following table of results shows that the concentrations for these analytes did not compare well.

Target Compound	Reported Concentration, ug/Kg	
	1 g Low Level	30 ul Medium Level Extract
acetone	3500	15000
2-butanone	5000	BDL
benzene	1200	5000
toluene	8600 *	110000
2-hexanone	12000 *	BDL
ethylbenzene	8000 *	100000
m,p-xylene	7400 *	170000
o-xylene	7900 *	110000
styrene	2300	23000

*estimated concentration of saturated compound
BDL - Below Detection Limit

We have concluded that these differences are due to a combination of sample inhomogeneity and the difference between the Low Level purge and trap and the Medium Level extraction procedure. In addition, because of the magnitude of the dilution, the 2-butanone and 2-hexanone are being reported only from the Low Level analysis, even though their concentrations are outside their analytical ranges. The data is being reported with reference to this qualifier.



Sarah A. Hubbard
Senior Volatile Data Specialist
January 3, 1990

QA Approval #1010
Linda Fowler
Employee #820
Sr. Quality Assurance Specialist
January 3, 1990

C. STANDARDS DATA

- (1) Initial Calibration Data (Form VI VOA) - in order by instrument if more than one instrument used.
 - (a) VOA standard(s) reconstructed ion chromatograms and quantitation reports (or legible facsimile) for the initial (five point) calibration. Spectra are not required.
 - (b) All initial calibration data must be included , regardless of when it was performed and for which case. When more than one initial calibration is performed , the data must be put in chronological order , by instrument.

- (2) Continuing Calibration (Form VII VOA) - in order by instrument , if more than one instrument used.
 - (a) VOA standard (s) reconstructed ion chromatograms and quantitation reports (or legible facsimile) for all continuing (12 hour) calibrations. Spectra are not required.
 - (b) When more than one continuing calibration is performed, forms must be in chronological order , within fraction and instrument.

- (3) Internal Standard Area Summary (Form VIII VOA) - in order by instrument , if more than one instrument used.

When more than one continuing calibration is performed , forms must be in chronological order , by instrument.

- (1) Initial Calibration Data (Form VI VOA) - in order by instrument if more than one instrument used.
 - (a) VOA standard(s) reconstructed ion chromatograms and quantitation reports (or legible facsimile) for the initial (five point) calibration. Spectra are not required.
 - (b) All initial calibration data must be included, regardless of when it was performed and for which case. When more than one initial calibration is performed, the data must be put in chronological order, by instrument.

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS
 Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Instrument ID: 13 Calibration Date(s): 12/28/89 12/28/89

Matrix: (soil/water) SOIL Level: (low/med) MED Column: (pack/cap) CAP

Min RRF for SPCC(#) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 30.0%

LAB FILE ID: RRF20 = CT891228A13 RRF50 = CV891228A13
 RRF100 = CW891228A13 RRF150 = CU891228A13 RRF200 = CS891228A13

COMPOUND	RRF20	RRF50	RRF100	RRF150	RRF200	RRF	% RSD
Chloromethane	0.922	0.670	0.512	0.672	0.871	0.729	22.9#
Bromomethane	1.421	1.124	0.895	1.062	1.106	1.122	17.0
Vinyl Chloride	1.079	0.833	0.646	0.869	1.059	0.897	19.9*
Chloroethane	0.688	0.574	0.465	0.536	0.556	0.564	14.3
Methylene Chloride	1.498	1.461	1.422	1.487	1.479	1.469	2.0
Acetone	0.205	0.274	0.250	0.278	0.232	0.248	12.2
Carbon Disulfide	3.675	4.009	3.209	2.528	1.513	2.987	33.3
1,1-Dichloroethene	1.232	1.266	1.087	0.733	0.619	0.987	29.9*
1,1-Dichloroethane	2.258	2.261	2.274	2.640	2.727	2.432	9.5#
1,2-Dichloroethene (total)	2.452	2.683	2.756	3.131	3.097	2.824	10.2
Chloroform	2.808	2.681	2.649	3.007	2.854	2.800	5.1*
1,2-Dichloroethane	1.831	1.779	1.830	2.110	2.128	1.936	8.7
2-Butanone	0.098	0.130	0.100	0.124	0.107	0.112	12.9
1,1,1-Trichloroethane	0.715	0.757	0.764	0.744	0.697	0.735	3.9
Carbon Tetrachloride	0.603	0.697	0.728	0.686	0.669	0.677	6.9
Vinyl Acetate	0.793	1.007	1.016	1.009	0.997	0.964	10.0
Bromodichloromethane	0.600	0.663	0.685	0.677	0.665	0.658	5.1
1,2-Dichloropropane	0.370	0.356	0.358	0.363	0.361	0.362	1.5*
cis-1,3-Dichloropropene	0.499	0.558	0.589	0.585	0.556	0.557	6.5
Trichloroethene	0.389	0.441	0.446	0.427	0.393	0.419	6.4
Dibromochloromethane	0.782	0.934	0.953	0.835	0.754	0.852	10.5
1,1,2-Trichloroethane	0.544	0.563	0.560	0.516	0.466	0.530	7.6
Benzene	0.822	0.886	0.848	0.878	0.858	0.858	3.0
Trans-1,3-Dichloropropene	0.412	0.475	0.462	0.453	0.500	0.460	7.0
Bromoform	0.603	0.806	0.894	0.750	0.655	0.742	15.7#
4-Methyl-2-Pentanone	0.179	0.192	0.207	0.235	0.217	0.206	10.6
2-Hexanone	0.207	0.244	0.217	0.227	0.217	0.222	6.3
Tetrachloroethene	0.487	0.559	0.548	0.532	0.480	0.521	6.9
1,1,2,2-Tetrachloroethane	0.653	0.694	0.682	0.684	0.705	0.684	2.8#
Toluene	0.255	0.278	0.294	0.329	0.320	0.295	10.3*
Chlorobenzene	0.862	0.983	0.951	0.959	0.968	0.945	5.0#
Ethylbenzene	0.391	0.414	0.414	0.408	0.400	0.405	2.4*
Styrene	0.950	1.121	1.008	1.029	0.990	1.020	6.2
Total Xylenes	1.166	1.376	1.238	1.291	1.251	1.264	6.1
Toluene-d8	0.461	0.495	0.501	0.552	0.543	0.510	7.3
BFB	0.898	0.936	0.873	0.879	0.892	0.896	2.8
1,2-Dichloroethane-d4	1.919	1.882	1.814	1.985	2.012	1.922	4.1

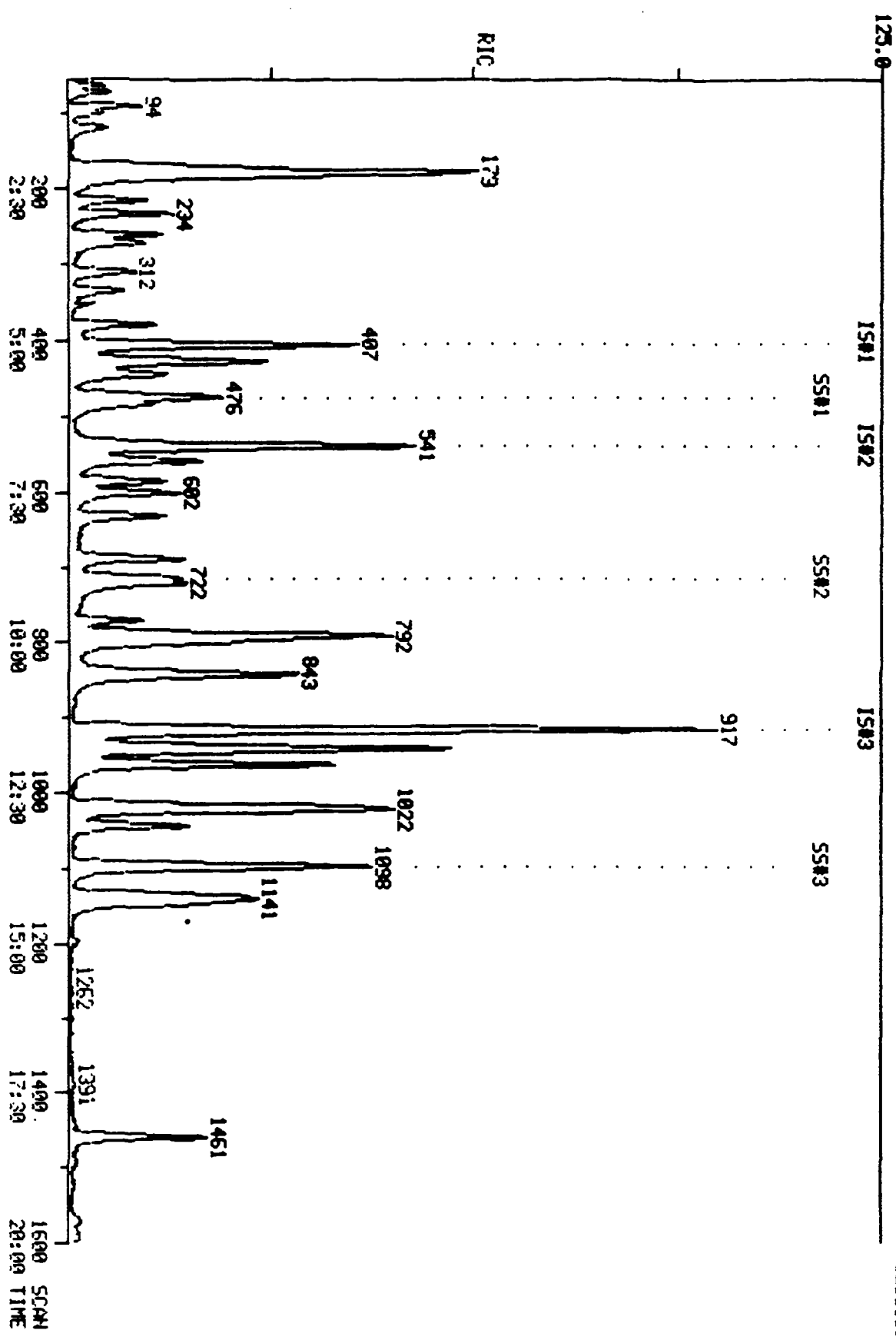
FORM VI VOA

1/87 Rev.

RIC
12/28/89 11:08:00
SAMPLE: SML USTD020 #1911 ON#13
CONDOS.:

COMPUchem LABS
COMPUchem DATA: CT891228A13 SCANS 57 TO 1600

120000.



QUANTITATION REPORT FILE: CT891228A13
DATA: CT891228A13.T1
12/28/89 11:08:00
SAMPLE: 5ML VSTD020 #1911 DN#13
CONDS. :
SUBMITTED BY: 13 ANALYST: 1492

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> NA#1
2	221 CHLOROMETHANE <74-87-3> NA#2
3	231 VINYL CHLORIDE <75-01-4> NA#3
4	220 BROMOMETHANE <78-83-9> NA#4
5	209 CHLOROETHANE <75-00-3> NA#5
6	230 TRICHLOROFLUOROMETHANE <75-69-4> NA#6
7	201 ACROLEIN <107-02-8> NA#7
8	216 1,1-DICHLOROETHENE <75-35-4> NA#8
9	254 CARBON DISULFIDE <75-15-0> NA#9
10	285 IODOMETHANE <74-88-4> NA#10
11	297 1,1,1-TRICHLORO-2,2,2-TRIFLUOROETHANE <354-38-5> NA#11
12	266 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE <76-13-1> NA#12
13	252 ACETONE (2-PROPANONE) <67-64-1> NA#13
14	*248 1,4-DIFLUOROBENZENE (IS) <540-36-3> NA#14
15	298 3-CHLOROPROPENE <107-05-1> NA#15
16	222 METHYLENE CHLORIDE <75-09-2> NA#16
17	226 TRANS-1,2-DICHLOROETHENE <156-60-5> NA#17
18	202 ACRYLONITRILE <107-13-1> NA#18
19	214 1,1-DICHLOROETHANE <75-34-3> NA#19
20	257 VINYL ACETATE <108-05-4> NA#20
21	237 CIS-1,2-DICHLOROETHENE <156-59-2> NA#21
22	253 2-BUTANONE <78-93-3> NA#22
23	211 CHLOROFORM <67-66-2> NA#23
24	227 1,1,1-TRICHLOROETHANE <71-55-6> NA#24
25	206 CARBON TETRACHLORIDE <56-23-5> NA#25
26	203 BENZENE <71-43-2> NA#26
27	215 1,2-DICHLOROETHANE <107-06-2> NA#27
28	272 CROTONALDEHYDE <4170-30-3> NA#28
29	*270 D5-CHLOROBENZENE (IS) <XX-XX-X> NA#29
30	229 TRICHLOROETHENE <79-01-6> NA#30
31	217 1,2-DICHLOROPROPANE <78-87-5> NA#31
32	286 DIBROMOMETHANE <74-95-3> NA#32
33	212 BROMODICHLOROMETHANE <75-27-4> NA#33
34	210 2-CHLOROETHYL VINYL ETHER <110-75-8> NA#34
35	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> NA#35
36	256 4-METHYL-2-PENTANONE <108-01-1> NA#36
37	225 TOLUENE <108-88-3> NA#37
38	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> NA#38
39	228 1,1,2-TRICHLOROETHANE <79-00-5> NA#39
40	287 ETHYLMETHACRYLATE <96-18-4> NA#40
41	224 TETRACHLOROETHENE <127-18-4> NA#41
42	255 2-HEXANONE <591-78-6> NA#42
43	208 DIBROMOCHLOROMETHANE <124-48-1> NA#43
44	245 1,2-DIBROMOETHANE <1060-93-4> NA#44
45	207 CHLOROBENZENE <108-90-7> NA#45
46	273 1,1,1,2-TETRACHLOROETHANE <630-20-6> NA#46

NO NAME
 47 219 ETHYLBENZENE <100-41-4> NA#47
 48 330 M,P-XYLENE <133-02-7> NA#48
 49 239 O-XYLENE <133-02-7> NA#49
 50 251 STYRENE <100-42-5> NA#50
 51 205 BROMOFORM <75-25-2> NA#51
 52 274 CIS-1,4-DICHLORO-2-BUTENE <764-71-0> NA#52
 53 275 1,2,3-TRICHLOROPROPANE <96-18-4> NA#53
 54 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> NA#54
 55 290 TRANS-1,4-DICHLORO-2-BUTENE <110-57-6> NA#55
 56 262 1,2-DIBROMO-3-CHLOROPROPANE <96-12-8> NA#56
 57 #258 D4-1,2-DICHLOROETHANE NA#57
 58 #247 BROMOFLUOROBENZENE <460-00-4> NA#58
 59 #233 D8-TOLUENE NA#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	407	5:05	1	1.000	A BB	46617.	50.000 UG/L	1.70
2	50	67	0:50	1	0.165	A BB	17194.	33.278 UG/L	1.13
3	62	76	0:57	1	0.187	A BB	20113.	41.206 UG/L	1.40
4	94	94	1:10	1	0.231	A BB	26492.	26.039 UG/L	0.88
5	64	103	1:17	1	0.253	A BB	12828.	25.521 UG/L	0.87
6	101	121	1:31	1	0.297	A BB	22113.	20.950 UG/L	0.71
7	56	173	2:10	1	0.425	A BV	17002.	545.937 UG/L	18.54
8	96	171	2:08	1	0.420	A BB	22967.	29.049 UG/L	0.99
9	76	179	2:14	1	0.440	A BB	68529.	87.581 UG/L	2.97
10	142	181	2:16	1	0.445	A BB	71895.	38.002 UG/L	1.29
11	117	178	2:13	1	0.437	A BB	22914.	28.658 UG/L	0.97
12	85	181	2:16	1	0.445	A BB	26119.	30.916 UG/L	1.05
13	43	196	2:27	1	0.482	M XX	3814.	27.835 UG/L	0.95
14	114	541	6:46	14	1.000	A BB	156951.	50.000 UG/L	1.70
15	76	217	2:43	1	0.533	A BB	11153.	36.904 UG/L	1.25
16	84	234	2:55	1	0.575	A BV	27937.	37.683 UG/L	1.28
17	96	261	3:16	1	0.641	A BB	22510.	19.066 UG/L	0.65
18	53	273	3:25	1	0.671	A BB	37642.	417.407 UG/L	14.17
19	63	312	3:54	1	0.767	A BB	42105.	21.888 UG/L	0.74
20	43	337	4:13	14	0.623	A BB	49760.	43.109 UG/L	1.46
21	96	381	4:46	1	0.936	A BB	23212.	15.368 UG/L	0.52
22	72	398	4:58	1	0.978	M XX	1820.	24.888 UG/L	0.85
23	83	428	5:21	1	1.052	A BB	52354.	21.314 UG/L	0.72
24	97	430	5:22	14	0.795	A BB	44893.	28.309 UG/L	0.96
25	117	446	5:34	14	0.824	A VB	37842.	25.121 UG/L	0.85
26	78	473	5:55	14	0.874	A BB	51590.	22.074 UG/L	0.75
27	62	486	6:04	1	1.194	A BB	34146.	22.892 UG/L	0.78
28	70	541	6:46	14	1.000	M XX	11842.	240.090 UG/L	8.15
29	117	916	11:27	29	1.000	A BB	215033.	50.000 UG/L	1.70
30	130	559	6:59	14	1.033	A BB	24440.	18.598 UG/L	0.63
31	63	587	7:20	14	1.085	A BB	23214.	25.808 UG/L	0.88
32	174	603	7:32	1	1.482	A BB	28633.	21.158 UG/L	0.72
33	83	634	7:55	14	1.172	A BB	37687.	19.976 UG/L	0.68
34	63	687	8:35	14	1.270	A BB	7812.	20.257 UG/L	0.69
35	75	691	8:38	14	1.277	A BB	31339.	16.186 UG/L	0.55
36	43	729	9:07	29	0.796	M XX	15409.	14.135 UG/L	0.48
37	92	723	9:02	29	0.789	A BB	21894.	7.724 UG/L	0.26
38	75	772	9:39	14	1.427	A BV	25875.	30.605 UG/L	1.04
39	97	795	9:56	14	1.470	A BB	34151.	29.571 UG/L	1.00
40	69	803	10:02	29	0.877	A BB	43242.	33.466 UG/L	1.14
41	164	790	9:52	29	0.862	A BB	41921.	19.403 UG/L	0.66

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA (HIGHT)	AMOUNT	%TOT
42	43	849	10:37	29	0.927	A BB	17781.	23.176 UG/L	0.79
43	129	842	10:31	14	1.556	A BB	49093.	28.643 UG/L	0.97
44	107	845	10:34	14	1.562	A BB	53655.	27.527 UG/L	0.93
45	112	920	11:30	29	1.004	A BB	74120.	17.743 UG/L	0.60
46	131	941	11:46	14	1.739	A BB	50447.	32.341 UG/L	1.10
47	106	945	11:49	29	1.032	A BV	33645.	18.786 UG/L	0.64
48	106	964	12:03	29	1.052	A VB	53353.	18.652 UG/L	0.63
49	106	1018	12:43	29	1.111	A BB	46934.	18.455 UG/L	0.63
50	104	1024	12:48	29	1.118	A BB	81705.	18.830 UG/L	0.64
51	173	1045	13:04	14	1.932	A BB	37826.	29.210 UG/L	0.99
52	88	1103	13:47	14	2.039	A BB	16719.	38.853 UG/L	1.32
53	110	1136	14:12	29	1.240	A BB	19282.	22.333 UG/L	0.76
54	83	1143	14:17	29	1.248	A BB	56168.	18.298 UG/L	0.62
55	53	1150	14:22	29	1.255	A BB	13551.	27.943 UG/L	0.95
56	157	1461	18:16	29	1.595	A BB	32001.	232.101 UG/L	7.88
57	65	478	5:58	1	1.174	A BB	35785.	26.495 UG/L	0.90
58	95	1098	13:43	29	1.199	A BB	77257.	19.766 UG/L	0.67
59	98	716	8:57	29	0.782	A BB	39622.	7.959 UG/L	0.27

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:08	0.99	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	0:50	1.00	10.000	0.02	33.28	50.00	0.369	0.554	0.67
3	0:56	1.01	10.000	0.02	41.21	50.00	0.431	0.524	0.82
4	1:11	0.99	10.000	0.02	26.04	50.00	0.568	1.091	0.52
5	1:17	1.00	10.000	0.03	25.52	50.00	0.275	0.539	0.51
6	1:34	0.96	10.000	0.03	20.95	50.00	0.474	1.132	0.42
7	2:16	0.95	90.000	0.00	545.94	500.01	0.036	0.033	1.09
8	2:10	0.98	5.000	0.08	29.05	50.00	0.493	0.848	0.58
9	2:17	0.98	5.000	0.09	87.58	50.00	1.470	0.839	1.75
10	2:17	0.99	10.000	0.04	38.00	50.00	1.542	2.029	0.76
11	2:13	1.00	10.000	0.04	28.66	50.00	0.492	0.858	0.57
12	2:15	1.01	10.000	0.04	30.92	50.00	0.560	0.906	0.62
13	2:45	0.89	10.000	0.05	27.83	50.00	0.082	0.147	0.56
14	6:46	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
15	2:44	0.99	10.000	0.05	36.90	50.00	0.239	0.324	0.74
16	2:59	0.98	5.000	0.11	37.68	50.00	0.599	0.795	0.75
17	3:18	0.99	5.000	0.13	19.07	50.00	0.483	1.266	0.38
18	3:35	0.95	120.000	0.01	417.41	500.01	0.081	0.097	0.83
19	3:56	0.99	5.000	0.15	21.89	50.00	0.903	2.063	0.44
20	4:17	0.98	10.000	0.06	43.11	50.00	0.317	0.368	0.86
21	4:49	0.99	5.000	0.19	15.37	50.00	0.498	1.620	0.31
22	5:16	0.95	10.000	0.10	24.89	50.00	0.039	0.078	0.50
23	5:25	0.99	5.000	0.21	21.31	50.00	1.123	2.635	0.43
24	5:24	1.00	5.000	0.16	28.31	50.00	0.286	0.505	0.57
25	5:35	1.00	5.000	0.16	25.12	50.00	0.241	0.480	0.50
26	5:57	0.99	5.000	0.17	22.07	50.00	0.329	0.745	0.44
27	6:07	0.99	5.000	0.24	22.89	50.00	0.732	1.600	0.46
28	6:58	0.97	100.000	0.01	240.09	500.01	0.008	0.016	0.48
29	11:26	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
30	7:01	1.00	5.000	0.21	18.60	50.00	0.156	0.419	0.37
31	7:23	0.99	5.000	0.22	25.81	50.00	0.148	0.287	0.52
32	7:34	1.00	5.000	0.30	21.16	50.00	0.614	1.451	0.42
33	7:57	1.00	5.000	0.23	19.98	50.00	0.240	0.601	0.40
34	8:37	1.00	10.000	0.13	20.26	50.00	0.050	0.123	0.41
35	8:40	1.00	5.000	0.26	16.19	50.00	0.200	0.617	0.32
36	9:11	0.99	15.000	0.05	14.14	50.00	0.072	0.253	0.28

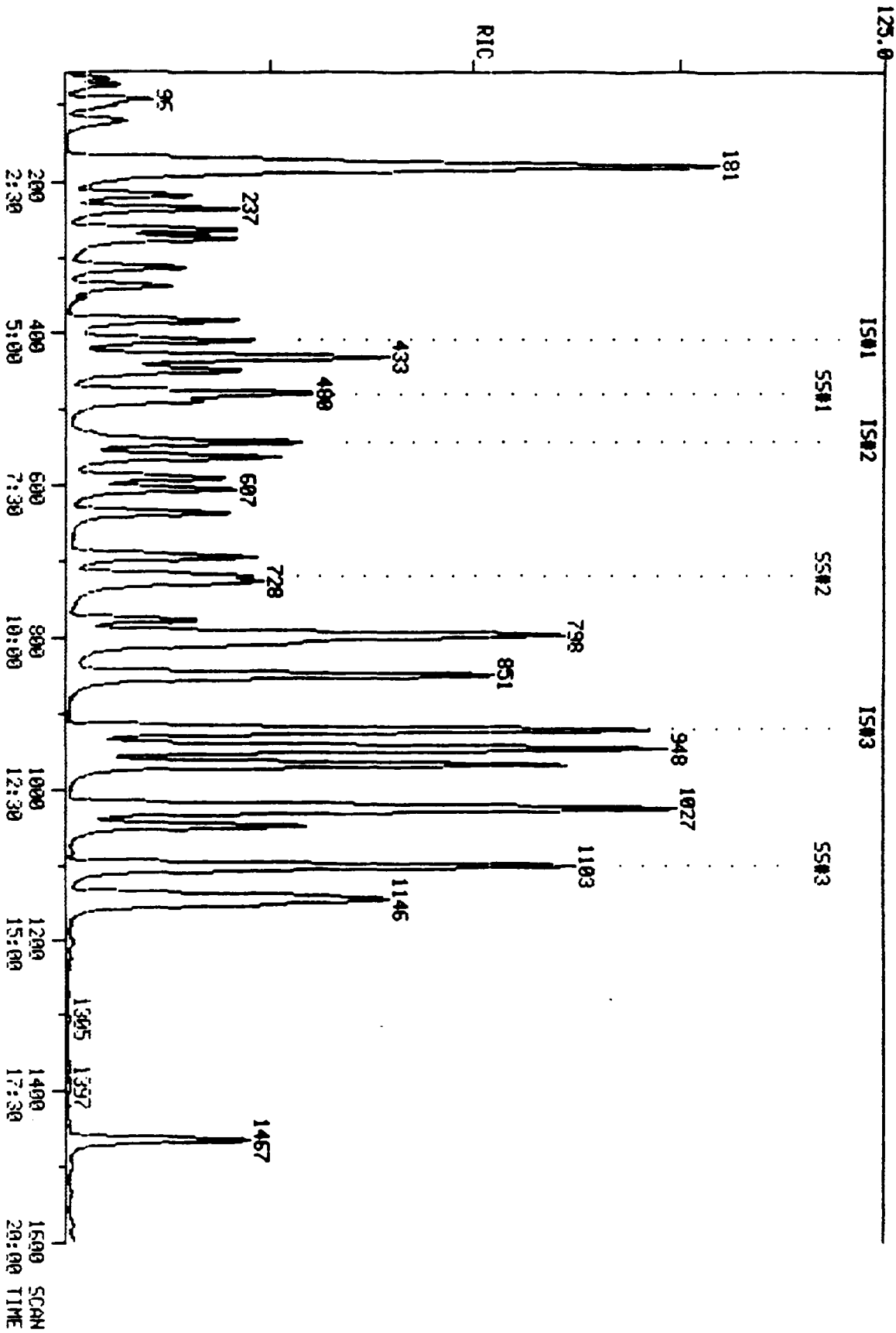
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
37	9:04	1.00	5.000	0.16	7.72	50.00	0.102	0.659	0.19
38	9:42	0.99	5.000	0.29	30.61	50.00	0.165	0.269	0.61
39	9:58	1.00	5.000	0.29	29.57	50.00	0.218	0.368	0.59
40	10:04	1.00	10.000	0.09	33.47	50.00	0.201	0.300	0.67
41	9:55	1.00	5.000	0.17	19.40	50.00	0.195	0.502	0.39
42	10:38	1.00	15.000	0.06	23.18	50.00	0.083	0.178	0.46
43	10:33	1.00	5.000	0.31	28.64	50.00	0.313	0.546	0.57
44	10:35	1.00	5.000	0.31	27.53	50.00	0.342	0.621	0.55
45	11:29	1.00	5.000	0.20	17.74	50.00	0.345	0.971	0.35
46	11:45	1.00	5.000	0.35	32.34	50.00	0.321	0.497	0.65
47	11:47	1.00	5.000	0.21	18.79	50.00	0.156	0.416	0.38
48	12:02	1.00	5.000	0.21	18.65	50.00	0.248	0.665	0.37
49	12:43	1.00	5.000	0.22	18.45	50.00	0.218	0.591	0.37
50	12:47	1.00	5.000	0.22	18.83	50.00	0.380	1.009	0.38
51	13:04	1.00	5.000	0.39	29.21	50.00	0.241	0.413	0.58
52	13:47	1.00	15.000	0.14	38.85	50.00	0.107	0.137	0.78
53	14:12	1.00	15.000	0.08	22.33	50.00	0.090	0.201	0.45
54	14:16	1.00	5.000	0.25	18.30	50.00	0.261	0.714	0.37
55	14:22	1.00	15.000	0.08	27.94	50.00	0.063	0.113	0.56
56	18:15	1.00	10.000	0.16	232.10	100.00	0.074	0.032	2.32
57	6:01	0.99	5.000	0.23	26.50	50.00	0.768	1.449	0.53
58	13:42	1.00	5.000	0.24	19.77	50.00	0.359	0.909	0.40
59	8:58	1.00	5.000	0.16	7.96	50.00	0.184	1.158	0.16

RIC
12/28/89 12:30:00
SAMPLE: SML UST0050 #1912 ON#13
CONDOS.:

COMPUchem LABS

COMPUchem DATA: CV891228A13 SCANS 59 TO 1600

191680.



QUANTITATION REPORT FILE: CV891228A13
DATA: CV891228A13.TI
12/28/89 12:30:00
SAMPLE: 5ML VSTD050 #1912 ON#13
CONDS.:
SUBMITTED BY: 13 ANALYST: 1492

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> NA#1
2	221 CHLOROMETHANE <74-87-3> NA#2
3	231 VINYL CHLORIDE <75-01-4> NA#3
4	220 BROMOMETHANE <78-83-9> NA#4
5	209 CHLOROETHANE <75-00-3> NA#5
6	230 TRICHLOROFLUOROMETHANE <75-69-4> NA#6
7	201 ACROLEIN <107-02-8> NA#7
8	216 1,1-DICHLOROETHENE <75-35-4> NA#8
9	254 CARBON DISULFIDE <75-15-0> NA#9
10	285 IODOMETHANE <74-88-4> NA#10
11	297 1,1,1-TRICHLORO-2,2,2-TRIFLUOROETHANE <354-58-5> NA#11
12	266 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE <76-13-1> NA#12
13	252 ACETONE (2-PROPANONE) <67-64-1> NA#13
14	*248 1,4-DIFLUOROBENZENE (IS) <340-36-3> NA#14
15	298 3-CHLOROPROPENE <107-09-1> NA#15
16	222 METHYLENE CHLORIDE <75-09-2> NA#16
17	226 TRANS-1,2-DICHLOROETHENE <156-60-5> NA#17
18	202 ACRYLONITRILE <107-13-1> NA#18
19	214 1,1-DICHLOROETHANE <75-34-3> NA#19
20	257 VINYL ACETATE <108-05-4> NA#20
21	237 CIS-1,2-DICHLOROETHENE <156-59-2> NA#21
22	253 2-BUTANONE <78-93-3> NA#22
23	211 CHLOROFORM <67-66-2> NA#23
24	227 1,1,1-TRICHLOROETHANE <71-55-6> NA#24
25	206 CARBON TETRACHLORIDE <56-23-5> NA#25
26	203 BENZENE <71-43-2> NA#26
27	215 1,2-DICHLOROETHANE <107-06-2> NA#27
28	272 CROTONALDEHYDE <4170-30-3> NA#28
29	*270 D5-CHLOROBENZENE (IS) <XX-XX-X> NA#29
30	229 TRICHLOROETHENE <79-01-6> NA#30
31	217 1,2-DICHLOROPROPANE <78-87-5> NA#31
32	286 DIBROMOMETHANE <74-95-3> NA#32
33	212 BROMODICHLOROMETHANE <75-27-4> NA#33
34	210 2-CHLOROETHYL VINYL ETHER <110-75-8> NA#34
35	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> NA#35
36	256 4-METHYL-2-PENTANONE <108-01-1> NA#36
37	225 TOLUENE <108-88-3> NA#37
38	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> NA#38
39	228 1,1,2-TRICHLOROETHANE <79-00-5> NA#39
40	287 ETHYLMETHACRYLATE <96-18-4> NA#40
41	224 TETRACHLOROETHENE <127-18-4> NA#41
42	255 2-HEXANONE <591-78-6> NA#42
43	208 DIBROMOCHLOROMETHANE <124-48-1> NA#43
44	245 1,2-DIBROMOETHANE <1060-93-4> NA#44
45	207 CHLOROBENZENE <108-90-7> NA#45
46	273 1,1,1,2-TETRACHLOROETHANE <630-20-6> NA#46

NO NAME
 47 219 ETHYLBENZENE <100-41-4> NA#47
 48 330 M,P-XYLENE <133-02-7> NA#48
 49 239 O-XYLENE <133-02-7> NA#49
 50 251 STYRENE <100-42-5> NA#50
 51 205 BROMOFORM <75-25-2> NA#51
 52 274 CIS-1,4-DICHLORO-2-BUTENE <764-71-0> NA#52
 53 275 1,2,3-TRICHLOROPROPANE <96-18-4> NA#53
 54 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> NA#54
 55 290 TRANS-1,4-DICHLORO-2-BUTENE <110-57-6> NA#55
 56 262 1,2-DIBROMO-3-CHLOROPROPANE <96-12-8> NA#56
 57 #258 D4-1,2-DICHLOROETHANE NA#57
 58 #247 BROMOFLUOROBENZENE <460-00-4> NA#58
 59 #233 DB-TOLUENE NA#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	411	5:08	1	1.000	A BB	49092.	50.000 UG/L	0.65
2	50	69	0:52	1	0.168	A BB	32881.	60.432 UG/L	0.79
3	62	78	0:58	1	0.190	A BB	40889.	79.546 UG/L	1.04
4	94	96	1:12	1	0.234	A BB	55164.	51.486 UG/L	0.67
5	64	104	1:18	1	0.253	A BB	28171.	53.221 UG/L	0.70
6	101	124	1:33	1	0.302	A BB	64061.	57.632 UG/L	0.75
7	56	176	2:12	1	0.428	A BB	50732.	1546.880 UG/L	20.25
8	96	173	2:10	1	0.421	A BB	62145.	74.639 UG/L	0.98
9	76	181	2:16	1	0.440	A BB	196827.	238.866 UG/L	3.13
10	142	183	2:17	1	0.445	A BB	171231.	85.945 UG/L	1.13
11	117	183	2:17	1	0.445	A BB	56453.	67.046 UG/L	0.88
12	85	183	2:17	1	0.445	A BB	57599.	64.741 UG/L	0.85
13	43	201	2:31	1	0.489	A BB	13442.	93.155 UG/L	1.22
14	114	544	6:48	14	1.000	A BB	155642.	50.000 UG/L	0.65
15	76	220	2:45	1	0.535	A BB	26751.	84.055 UG/L	1.10
16	84	237	2:58	1	0.577	A BB	71711.	91.850 UG/L	1.20
17	96	264	3:18	1	0.642	A BB	63194.	50.827 UG/L	0.67
18	53	277	3:28	1	0.674	A BB	126119.	1327.990 UG/L	17.39
19	63	315	3:56	1	0.766	A BB	110996.	54.792 UG/L	0.72
20	43	339	4:14	14	0.623	A BB	156787.	136.971 UG/L	1.79
21	96	384	4:48	1	0.934	A BB	68551.	43.098 UG/L	0.56
22	72	404	5:03	1	0.983	A BB	6403.	83.145 UG/L	1.09
23	83	432	5:24	1	1.051	A BB	131627.	50.886 UG/L	0.67
24	97	434	5:25	14	0.798	A BB	117834.	74.929 UG/L	0.98
25	117	450	5:37	14	0.827	A VB	108525.	72.647 UG/L	0.95
26	78	478	5:58	14	0.879	A BB	137824.	59.466 UG/L	0.78
27	62	490	6:07	1	1.192	A BB	87350.	55.608 UG/L	0.73
28	70	542	6:46	14	0.996	A BB	27326.	558.688 UG/L	7.32
29	117	921	11:31	29	1.000	A BB	214975.	50.000 UG/L	0.65
30	130	564	7:03	14	1.037	A BB	68586.	52.632 UG/L	0.69
31	63	592	7:24	14	1.088	A BB	55390.	62.098 UG/L	0.81
32	174	607	7:35	1	1.477	A BB	69389.	48.689 UG/L	0.64
33	83	638	7:58	14	1.173	A BB	103207.	55.163 UG/L	0.72
34	63	692	8:39	14	1.272	A BB	21721.	56.797 UG/L	0.74
35	75	696	8:42	14	1.279	A BB	86863.	45.241 UG/L	0.59
36	43	732	9:09	29	0.795	A BB	41181.	37.788 UG/L	0.49
37	92	728	9:06	29	0.790	A BB	59721.	21.075 UG/L	0.28
38	75	778	9:43	14	1.430	A BB	73893.	88.137 UG/L	1.15
39	97	800	10:00	14	1.471	A BB	87554.	76.450 UG/L	1.00
40	69	809	10:07	29	0.878	A BB	106009.	82.065 UG/L	1.07
41	164	796	9:57	29	0.864	A BB	120239.	55.667 UG/L	0.73

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
42	43	854	10:40	29	0.927	A BB	52500.	68.448 UG/L	0.90
43	129	849	10:37	14	1.561	A BB	145428.	85.562 UG/L	1.12
44	107	852	10:39	14	1.566	A BB	135624.	70.164 UG/L	0.92
45	112	925	11:34	29	1.004	A BB	211271.	50.588 UG/L	0.66
46	131	946	11:49	14	1.739	A BB	125021.	80.825 UG/L	1.06
47	106	949	11:52	29	1.030	A BV	89104.	49.767 UG/L	0.65
48	106	969	12:07	29	1.052	A VB	158362.	55.379 UG/L	0.73
49	106	1023	12:47	29	1.111	A BB	137471.	54.069 UG/L	0.71
50	104	1029	12:52	29	1.117	A BB	241045.	55.568 UG/L	0.73
51	173	1050	13:07	14	1.930	A BB	125477.	97.710 UG/L	1.28
52	88	1107	13:50	14	2.035	A BB	41412.	97.047 UG/L	1.27
53	110	1141	14:16	29	1.239	A BB	45674.	52.916 UG/L	0.69
54	83	1148	14:21	29	1.246	A BB	149173.	48.608 UG/L	0.64
55	53	1155	14:26	29	1.254	A BB	33741.	69.595 UG/L	0.91
56	157	1467	18:20	29	1.593	A BB	70698.	512.907 UG/L	6.72
57	65	482	6:01	1	1.173	A BB	92403.	64.966 UG/L	0.85
58	95	1102	13:46	29	1.197	A BB	201164.	51.480 UG/L	0.67
59	98	721	9:01	29	0.783	A BB	106455.	21.390 UG/L	0.28

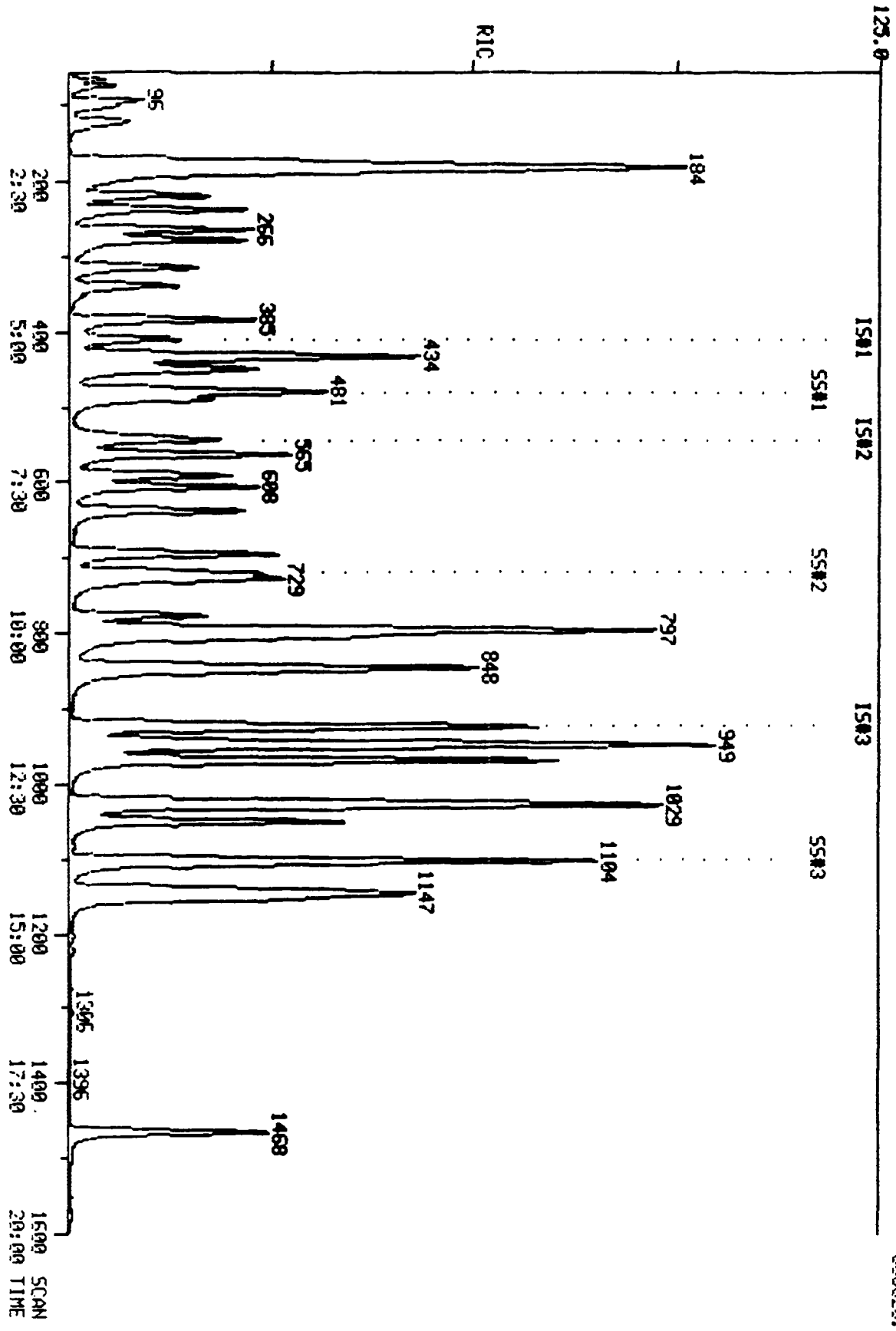
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:08	1.00	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	0:50	1.03	10.000	0.02	60.43	50.00	0.670	0.554	1.21
3	0:56	1.04	10.000	0.02	79.55	50.00	0.833	0.524	1.59
4	1:11	1.01	10.000	0.02	51.49	50.00	1.124	1.091	1.03
5	1:17	1.01	10.000	0.03	53.22	50.00	0.574	0.539	1.06
6	1:34	0.98	10.000	0.03	57.63	50.00	1.305	1.132	1.15
7	2:16	0.97	90.000	0.00	1546.89	500.01	0.103	0.033	3.09
8	2:10	0.99	5.000	0.08	74.64	50.00	1.266	0.848	1.49
9	2:17	0.99	5.000	0.09	238.87	50.00	4.009	0.839	4.78
10	2:17	1.00	10.000	0.04	85.95	50.00	3.488	2.029	1.72
11	2:13	1.03	10.000	0.04	67.05	50.00	1.150	0.858	1.34
12	2:15	1.02	10.000	0.04	64.74	50.00	1.173	0.906	1.29
13	2:45	0.91	10.000	0.05	93.15	50.00	0.274	0.147	1.86
14	6:46	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
15	2:44	1.00	10.000	0.05	84.05	50.00	0.545	0.324	1.68
16	2:59	0.99	5.000	0.12	91.85	50.00	1.461	0.795	1.84
17	3:18	1.00	5.000	0.13	50.83	50.00	1.287	1.266	1.02
18	3:35	0.97	120.000	0.01	1328.00	500.01	0.257	0.097	2.66
19	3:56	1.00	5.000	0.15	54.79	50.00	2.261	2.063	1.10
20	4:17	0.99	10.000	0.06	136.97	50.00	1.007	0.368	2.74
21	4:49	1.00	5.000	0.19	43.10	50.00	1.396	1.620	0.86
22	5:16	0.96	10.000	0.10	83.15	50.00	0.130	0.078	1.66
23	5:25	1.00	5.000	0.21	50.89	50.00	2.681	2.635	1.02
24	5:24	1.00	5.000	0.16	74.93	50.00	0.757	0.505	1.50
25	5:35	1.01	5.000	0.17	72.65	50.00	0.697	0.480	1.45
26	5:57	1.00	5.000	0.18	59.47	50.00	0.886	0.745	1.19
27	6:07	1.00	5.000	0.24	55.61	50.00	1.779	1.600	1.11
28	6:58	0.97	100.000	0.01	558.69	500.01	0.018	0.016	1.12
29	11:26	1.01	5.000	0.20	50.00	50.00	1.000	1.000	1.00
30	7:01	1.01	5.000	0.21	52.63	50.00	0.441	0.419	1.05
31	7:23	1.00	5.000	0.22	62.10	50.00	0.356	0.287	1.24
32	7:34	1.00	5.000	0.30	48.69	50.00	1.413	1.451	0.97
33	7:57	1.00	5.000	0.23	55.16	50.00	0.663	0.601	1.10
34	8:37	1.00	10.000	0.13	56.80	50.00	0.140	0.123	1.14
35	8:40	1.00	5.000	0.26	45.24	50.00	0.558	0.617	0.90
36	9:11	1.00	15.000	0.05	37.79	50.00	0.192	0.253	0.76

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
37	9:04	1.00	5.000	0.16	21.07	50.00	0.278	0.659	0.42
38	9:42	1.00	5.000	0.29	88.14	50.00	0.475	0.269	1.76
39	9:58	1.00	5.000	0.29	76.45	50.00	0.563	0.368	1.53
40	10:04	1.00	10.000	0.09	82.07	50.00	0.493	0.300	1.64
41	9:55	1.00	5.000	0.17	55.67	50.00	0.559	0.502	1.11
42	10:38	1.00	15.000	0.06	68.45	50.00	0.244	0.178	1.37
43	10:33	1.01	5.000	0.31	85.56	50.00	0.934	0.546	1.71
44	10:35	1.01	5.000	0.31	70.16	50.00	0.871	0.621	1.40
45	11:29	1.01	5.000	0.20	50.59	50.00	0.983	0.971	1.01
46	11:45	1.01	5.000	0.35	80.82	50.00	0.803	0.497	1.62
47	11:47	1.01	5.000	0.21	49.77	50.00	0.414	0.416	1.00
48	12:02	1.01	5.000	0.21	55.38	50.00	0.737	0.665	1.11
49	12:43	1.00	5.000	0.22	54.07	50.00	0.639	0.591	1.08
50	12:47	1.01	5.000	0.22	55.57	50.00	1.121	1.009	1.11
51	13:04	1.00	5.000	0.39	97.71	50.00	0.806	0.413	1.95
52	13:47	1.00	15.000	0.14	97.05	50.00	0.266	0.137	1.94
53	14:12	1.00	15.000	0.08	52.92	50.00	0.212	0.201	1.06
54	14:16	1.01	5.000	0.25	48.61	50.00	0.694	0.714	0.97
55	14:22	1.00	15.000	0.08	69.60	50.00	0.157	0.113	1.39
56	18:15	1.00	10.000	0.16	512.91	100.00	0.164	0.032	5.13
57	6:01	1.00	5.000	0.23	64.97	50.00	1.882	1.449	1.30
58	13:42	1.01	5.000	0.24	51.48	50.00	0.936	0.909	1.03
59	8:58	1.01	5.000	0.16	21.39	50.00	0.495	1.158	0.43

RIC
 12/28/89 13:04:00
 SAMPLE: SML USTD100 #1913 ON#13
 COND.S.:

COMPUCHEM LABS
 COMPUCHEM DATA: CM891228A13 SCANS 58 TO 1600

363520.



QUANTITATION REPORT FILE: CW891228A13
DATA: CW891228A13.TI
12/28/89 13:04:00
SAMPLE: 5ML VSTD100 #1913 ON#13
CONDS.:
SUBMITTED BY: 13 ANALYST: 1492

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> NA#1
2	221 CHLOROMETHANE <74-87-3> NA#2
3	231 VINYL CHLORIDE <75-01-4> NA#3
4	220 BROMOMETHANE <78-83-9> NA#4
5	209 CHLOROETHANE <75-00-3> NA#5
6	230 TRICHLOROFLUOROMETHANE <75-69-4> NA#6
7	201 ACROLEIN <107-02-8> NA#7
8	216 1,1-DICHLOROETHENE <75-35-4> NA#8
9	254 CARBON DIBULFIDE <75-15-0> NA#9
10	285 IODOMETHANE <74-88-4> NA#10
11	297 1,1,1-TRICHLORO-2,2,2-TRIFLUOROETHANE <354-58-5> NA#11
12	266 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE <76-13-1> NA#12
13	252 ACETONE (2-PROPANONE) <67-64-1> NA#13
14	*248 1,4-DIFLUOROBENZENE (IS) <340-36-3> NA#14
15	298 3-CHLOROPROPENE <107-05-1> NA#15
16	222 METHYLENE CHLORIDE <75-09-2> NA#16
17	226 TRANS-1,2-DICHLOROETHENE <156-60-5> NA#17
18	202 ACRYLONITRILE <107-13-1> NA#18
19	214 1,1-DICHLOROETHANE <75-34-3> NA#19
20	257 VINYL ACETATE <108-05-4> NA#20
21	237 CIS-1,2-DICHLOROETHENE <156-59-2> NA#21
22	253 2-BUTANONE <78-93-3> NA#22
23	211 CHLOROFORM <67-66-2> NA#23
24	227 1,1,1-TRICHLOROETHANE <71-55-6> NA#24
25	206 CARBON TETRACHLORIDE <56-23-5> NA#25
26	203 BENZENE <71-43-2> NA#26
27	215 1,2-DICHLOROETHANE <107-06-2> NA#27
28	272 CROTONALDEHYDE <4170-30-3> NA#28
29	*270 D5-CHLOROENZENE (IS) <XX-XX-X> NA#29
30	229 TRICHLOROETHENE <79-01-6> NA#30
31	217 1,2-DICHLOROPROPANE <78-87-5> NA#31
32	286 DIBROMOMETHANE <74-95-3> NA#32
33	212 BROMODICHLOROMETHANE <75-27-4> NA#33
34	210 2-CHLOROETHYL VINYL ETHER <110-75-8> NA#34
35	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> NA#35
36	256 4-METHYL-2-PENTANONE <108-01-1> NA#36
37	225 TOLUENE <108-88-3> NA#37
38	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> NA#38
39	228 1,1,2-TRICHLOROETHANE <79-00-5> NA#39
40	287 ETHYLMETHACRYLATE <96-18-4> NA#40
41	224 TETRACHLOROETHENE <127-18-4> NA#41
42	255 2-HEXANONE <591-78-6> NA#42
43	208 DIBROMOCHLOROMETHANE <124-48-1> NA#43
44	245 1,2-DIBROMOETHANE <1060-93-4> NA#44
45	207 CHLOROBENZENE <108-90-7> NA#45
46	273 1,1,1,2-TETRACHLOROETHANE <630-20-6> NA#46

NO NAME
 47 219 ETHYLBENZENE <100-41-4> NA#47
 48 330 M,P-XYLENE <133-02-7> NA#48
 49 239 O-XYLENE <133-02-7> NA#49
 50 251 STYRENE <100-42-5> NA#50
 51 205 BROMOFORM <75-25-2> NA#51
 52 274 CIS-1,4-DICHLORO-2-BUTENE <764-71-0> NA#52
 53 275 1,2,3-TRICHLOROPROPANE <96-18-4> NA#53
 54 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> NA#54
 55 290 TRANS-1,4-DICHLORO-2-BUTENE <110-57-6> NA#55
 56 262 1,2-DIBROMO-3-CHLOROPROPANE <96-12-8> NA#56
 57 #258 D4-1,2-DICHLOROETHANE NA#57
 58 #247 BROMOFLUOROBENZENE <460-00-4> NA#58
 59 #233 D8-TOLUENE NA#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HOHT)	AMOUNT	%TOT
1	128	411	5:08	1	1.000	A BB	50862.	50.000 UG/L	0.36
2	50	68	0:51	1	0.165	A BB	52040.	92.316 UG/L	0.66
3	62	77	0:58	1	0.187	A BB	65721.	123.406 UG/L	0.88
4	94	96	1:12	1	0.234	A BB	90994.	81.972 UG/L	0.58
5	64	104	1:18	1	0.253	A BB	47273.	86.201 UG/L	0.61
6	101	124	1:33	1	0.302	A BB	121361.	105.382 UG/L	0.75
7	56	178	2:13	1	0.433	A BB	88287.	2598.300 UG/L	18.48
8	96	174	2:10	1	0.423	A BB	110591.	128.203 UG/L	0.91
9	76	182	2:16	1	0.443	A BB	326410.	382.341 UG/L	2.72
10	142	184	2:18	1	0.448	A BB	328150.	158.975 UG/L	1.13
11	117	184	2:18	1	0.448	A BB	104825.	120.162 UG/L	0.85
12	85	184	2:18	1	0.448	A BB	113353.	122.975 UG/L	0.87
13	43	206	2:34	1	0.501	A BB	25393.	169.856 UG/L	1.21
14	114	546	6:49	14	1.000	A BB	159946.	50.000 UG/L	0.36
15	76	220	2:45	1	0.535	A BB	59358.	180.017 UG/L	1.28
16	84	239	2:59	1	0.582	A BB	144666.	178.847 UG/L	1.27
17	96	266	3:19	1	0.647	A BB	132919.	103.186 UG/L	0.73
18	53	280	3:30	1	0.681	A BB	247761.	2518.050 UG/L	17.91
19	63	316	3:57	1	0.769	A BB	231318.	110.213 UG/L	0.78
20	43	341	4:16	14	0.625	A BB	325034.	276.313 UG/L	1.97
21	96	385	4:49	1	0.937	A BB	147368.	89.427 UG/L	0.64
22	72	405	5:04	1	0.985	A BB	10127.	126.926 UG/L	0.90
23	83	434	5:25	1	1.056	A BB	269496.	100.560 UG/L	0.72
24	97	434	5:25	14	0.795	A BB	244479.	151.277 UG/L	1.08
25	117	451	5:38	14	0.826	A VB	232950.	151.741 UG/L	1.08
26	78	479	5:59	14	0.877	A BB	271170.	113.852 UG/L	0.81
27	62	492	6:09	1	1.197	A BB	186189.	114.405 UG/L	0.81
28	70	543	6:47	14	0.995	A BB	48189.	958.710 UG/L	6.82
29	117	922	11:31	29	1.000	A BB	223864.	50.000 UG/L	0.36
30	130	565	7:04	14	1.035	A BB	142743.	106.591 UG/L	0.76
31	63	593	7:25	14	1.086	A BB	114512.	124.924 UG/L	0.89
32	174	608	7:36	1	1.479	A BB	150879.	102.186 UG/L	0.73
33	83	640	8:00	14	1.172	A BB	219106.	113.959 UG/L	0.81
34	63	692	8:39	14	1.267	A BB	46321.	117.863 UG/L	0.84
35	75	697	8:43	14	1.277	A BB	188559.	95.565 UG/L	0.68
36	43	733	9:10	29	0.795	A BB	92619.	81.612 UG/L	0.58
37	92	729	9:07	29	0.791	A BB	131497.	44.561 UG/L	0.32
38	75	778	9:43	14	1.425	A BB	147905.	171.669 UG/L	1.22
39	97	800	10:00	14	1.465	A BB	179186.	152.250 UG/L	1.08
40	69	808	10:06	29	0.876	A BB	229941.	170.938 UG/L	1.22
41	164	796	9:57	29	0.863	A BB	245455.	109.126 UG/L	0.78

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
42	43	853	10:40	29	0.925	A BB	96936.	121.364 UG/L	0.86
43	129	846	10:34	14	1.549	A BB	304793.	174.498 UG/L	1.24
44	107	849	10:37	14	1.555	A BB	275261.	138.572 UG/L	0.99
45	112	926	11:34	29	1.004	A BB	425771.	97.901 UG/L	0.70
46	131	946	11:49	14	1.733	A BB	258819.	162.822 UG/L	1.16
47	106	951	11:53	29	1.031	A BV	185426.	99.452 UG/L	0.71
48	106	970	12:07	29	1.052	A VB	298588.	100.269 UG/L	0.71
49	106	1029	12:49	29	1.112	A BB	255651.	96.559 UG/L	0.69
50	104	1031	12:53	29	1.118	A BB	451242.	99.894 UG/L	0.71
51	173	1051	13:08	14	1.925	A BB	286025.	216.736 UG/L	1.54
52	88	1109	13:52	14	2.031	A BB	87342.	199.174 UG/L	1.42
53	110	1143	14:17	29	1.240	A BB	97398.	108.361 UG/L	0.77
54	83	1149	14:22	29	1.246	A BB	305238.	95.513 UG/L	0.68
55	53	1156	14:27	29	1.254	A BB	67438.	133.576 UG/L	0.95
56	157	1467	18:20	29	1.591	A BB	152773.	1064.340 UG/L	7.57
57	65	483	6:02	1	1.175	A BB	184564.	125.245 UG/L	0.89
58	95	1103	13:47	29	1.196	A BB	390703.	96.015 UG/L	0.68
59	98	721	9:01	29	0.782	A BB	224449.	43.307 UG/L	0.31

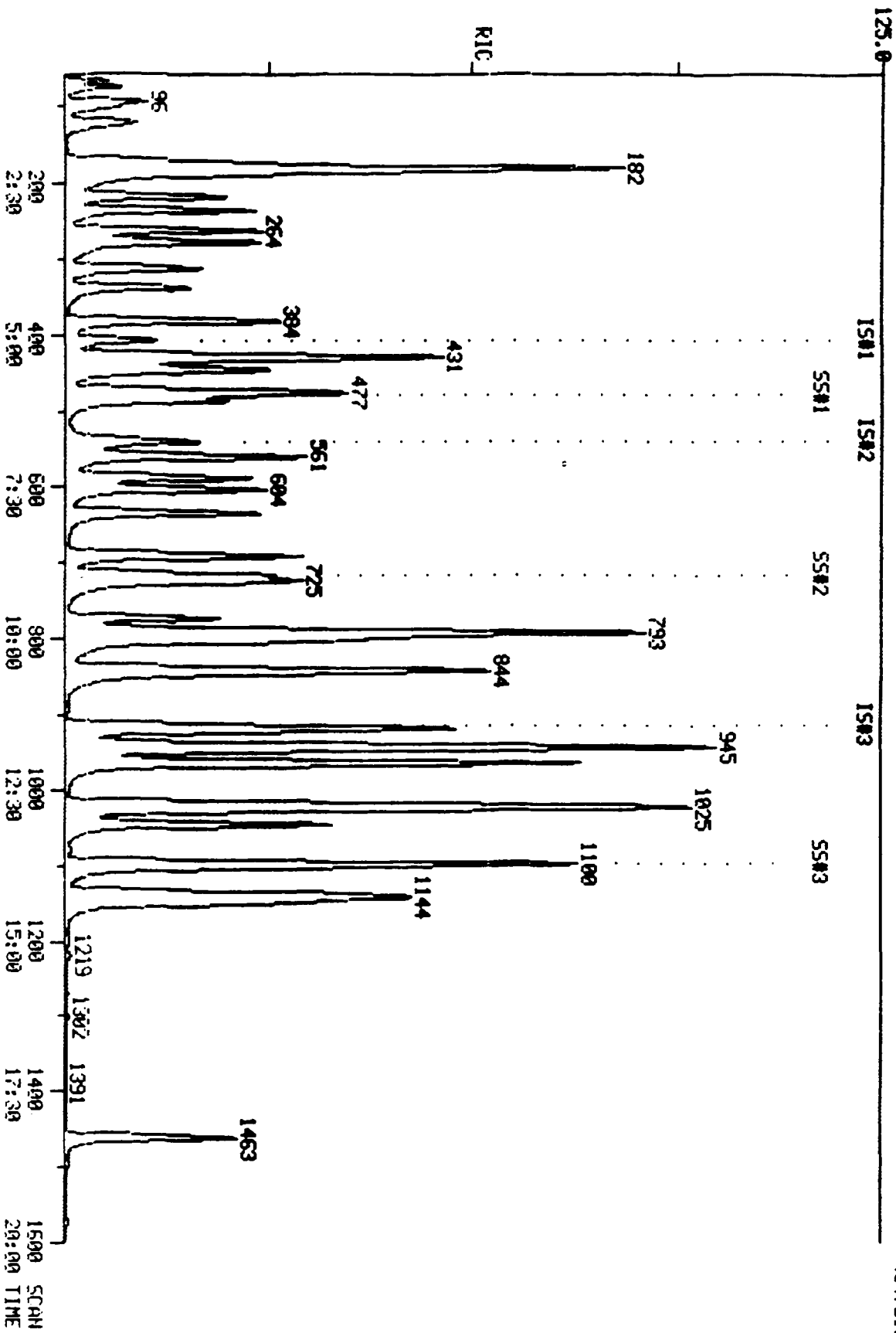
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:08	1.00	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	0:50	1.01	10.000	0.02	92.32	50.00	1.023	0.554	1.85
3	0:56	1.03	10.000	0.02	123.41	50.00	1.292	0.524	2.47
4	1:11	1.01	10.000	0.02	81.97	50.00	1.789	1.091	1.64
5	1:17	1.01	10.000	0.03	86.20	50.00	0.929	0.539	1.72
6	1:34	0.98	10.000	0.03	105.38	50.00	2.386	1.132	2.11
7	2:16	0.98	90.000	0.00	2598.31	500.01	0.174	0.033	5.20
8	2:10	1.00	5.000	0.08	128.20	50.00	2.174	0.848	2.56
9	2:17	0.99	5.000	0.09	382.34	50.00	6.418	0.839	7.65
10	2:17	1.01	10.000	0.04	158.98	50.00	6.452	2.029	3.18
11	2:13	1.03	10.000	0.04	120.16	50.00	2.061	0.858	2.40
12	2:15	1.02	10.000	0.04	122.97	50.00	2.229	0.906	2.46
13	2:45	0.94	10.000	0.05	169.86	50.00	0.499	0.147	3.40
14	6:46	1.01	5.000	0.20	50.00	50.00	1.000	1.000	1.00
15	2:44	1.00	10.000	0.05	180.02	50.00	1.167	0.324	3.60
16	2:59	1.00	5.000	0.12	178.85	50.00	2.844	0.795	3.58
17	3:18	1.01	5.000	0.13	103.19	50.00	2.613	1.266	2.06
18	3:35	0.98	120.000	0.01	2518.06	500.01	0.487	0.097	5.04
19	3:56	1.00	5.000	0.15	110.21	50.00	4.548	2.063	2.20
20	4:17	0.99	10.000	0.06	276.31	50.00	2.032	0.368	5.53
21	4:49	1.00	5.000	0.19	89.43	50.00	2.897	1.620	1.79
22	5:16	0.96	10.000	0.10	126.93	50.00	0.199	0.078	2.54
23	5:25	1.00	5.000	0.21	100.56	50.00	5.299	2.635	2.01
24	5:24	1.00	5.000	0.16	151.28	50.00	1.529	0.505	3.03
25	5:35	1.01	5.000	0.17	151.74	50.00	1.456	0.480	3.03
26	5:57	1.01	5.000	0.18	113.85	50.00	1.695	0.745	2.28
27	6:07	1.00	5.000	0.24	114.41	50.00	3.661	1.600	2.29
28	6:58	0.97	100.000	0.01	958.71	500.01	0.030	0.016	1.92
29	11:26	1.01	5.000	0.20	50.00	50.00	1.000	1.000	1.00
30	7:01	1.01	5.000	0.21	106.59	50.00	0.892	0.419	2.13
31	7:23	1.00	5.000	0.22	124.92	50.00	0.716	0.287	2.50
32	7:34	1.00	5.000	0.30	102.19	50.00	2.966	1.451	2.04
33	7:57	1.01	5.000	0.23	113.96	50.00	1.370	0.601	2.28
34	8:37	1.00	10.000	0.13	117.86	50.00	0.290	0.123	2.36
35	8:40	1.01	5.000	0.26	95.56	50.00	1.179	0.617	1.91
36	9:11	1.00	15.000	0.05	81.61	50.00	0.414	0.253	1.63

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
37	9:04	1.00	5.000	0.16	44.56	50.00	0.587	0.659	0.89
38	9:42	1.00	5.000	0.28	171.67	50.00	0.925	0.269	3.43
39	9:58	1.00	5.000	0.29	152.25	50.00	1.120	0.368	3.04
40	10:04	1.00	10.000	0.09	170.94	50.00	1.027	0.300	3.42
41	9:55	1.00	5.000	0.17	109.13	50.00	1.096	0.502	2.18
42	10:38	1.00	15.000	0.06	121.36	50.00	0.433	0.178	2.43
43	10:33	1.00	5.000	0.31	174.50	50.00	1.906	0.546	3.49
44	10:35	1.00	5.000	0.31	138.57	50.00	1.721	0.621	2.77
45	11:29	1.01	5.000	0.20	97.90	50.00	1.902	0.971	1.96
46	11:45	1.01	5.000	0.35	162.82	50.00	1.618	0.497	3.26
47	11:47	1.01	5.000	0.21	99.45	50.00	0.828	0.416	1.99
48	12:02	1.01	5.000	0.21	100.27	50.00	1.334	0.665	2.01
49	12:43	1.01	5.000	0.22	96.56	50.00	1.142	0.591	1.93
50	12:47	1.01	5.000	0.22	99.89	50.00	2.016	1.009	2.00
51	13:04	1.01	5.000	0.38	216.74	50.00	1.788	0.413	4.33
52	13:47	1.01	15.000	0.14	199.17	50.00	0.546	0.137	3.98
53	14:12	1.01	15.000	0.08	108.36	50.00	0.435	0.201	2.17
54	14:16	1.01	5.000	0.25	95.51	50.00	1.363	0.714	1.91
55	14:22	1.01	15.000	0.08	133.58	50.00	0.301	0.113	2.67
56	18:15	1.00	10.000	0.16	1064.35	100.00	0.341	0.032	10.64
57	6:01	1.00	5.000	0.24	125.25	50.00	3.629	1.449	2.50
58	13:42	1.01	5.000	0.24	96.02	50.00	1.745	0.909	1.92
59	8:58	1.01	5.000	0.16	43.31	50.00	1.003	1.158	0.87

RIC
12/28/89 11:52:00
SAMPLE: SML U5TD150 #1914 ON#13
CONDOS.:

COMPUCHEN LABS
COMPUCHEN DATA: CUB91228A13 SCANS 58 TO 1600

494729.



QUANTITATION REPORT FILE: CUB91228A13
DATA: CUB91228A13.TI
12/28/89 11:52:00
SAMPLE: 5ML VSTD150 #1914 ON#13
CONDS. :
SUBMITTED BY: 13 ANALYST: 1492

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> NA#1
2	221 CHLOROMETHANE <74-87-3> NA#2
3	231 VINYL CHLORIDE <75-01-4> NA#3
4	220 BROMOMETHANE <78-83-9> NA#4
5	209 CHLOROETHANE <75-00-3> NA#5
6	230 TRICHLOROFLUOROMETHANE <75-69-4> NA#6
7	201 ACROLEIN <107-02-8> NA#7
8	216 1,1-DICHLOROETHENE <75-35-4> NA#8
9	254 CARBON DISULFIDE <75-15-0> NA#9
10	285 IODOMETHANE <74-88-4> NA#10
11	297 1,1,1-TRICHLORO-2,2,2-TRIFLUOROETHANE <354-58-5> NA#11
12	266 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE <76-13-1> NA#12
13	252 ACETONE (2-PROPANONE) <67-64-1> NA#13
14	*248 1,4-DIFLUOROBENZENE (IS) <540-36-3> NA#14
15	298 3-CHLOROPROPENE <107-05-1> NA#15
16	222 METHYLENE CHLORIDE <75-09-2> NA#16
17	226 TRANS-1,2-DICHLOROETHENE <156-60-5> NA#17
18	202 ACRYLONITRILE <107-13-1> NA#18
19	214 1,1-DICHLOROETHANE <75-34-3> NA#19
20	257 VINYL ACETATE <108-05-4> NA#20
21	237 CIS-1,2-DICHLOROETHENE <156-59-2> NA#21
22	253 2-BUTANONE <78-93-3> NA#22
23	211 CHLOROFORM <67-66-2> NA#23
24	227 1,1,1-TRICHLOROETHANE <71-55-6> NA#24
25	206 CARBON TETRACHLORIDE <56-23-5> NA#25
26	203 BENZENE <71-43-2> NA#26
27	215 1,2-DICHLOROETHANE <107-06-2> NA#27
28	272 CROTONALDEHYDE <4170-30-3> NA#28
29	*270 D5-CHLOROBENZENE (IS) <XX-XX-X> NA#29
30	229 TRICHLOROETHENE <79-01-6> NA#30
31	217 1,2-DICHLOROPROPANE <78-87-5> NA#31
32	286 DIBROMOMETHANE <74-95-3> NA#32
33	212 BROMODICHLOROMETHANE <75-27-4> NA#33
34	210 2-CHLOROETHYL VINYL ETHER <110-75-8> NA#34
35	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> NA#35
36	256 4-METHYL-2-PENTANONE <108-01-1> NA#36
37	225 TOLUENE <108-88-3> NA#37
38	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> NA#38
39	228 1,1,2-TRICHLOROETHANE <79-00-5> NA#39
40	287 ETHYLMETHACRYLATE <96-18-4> NA#40
41	224 TETRACHLOROETHENE <127-18-4> NA#41
42	255 2-HEXANONE <591-78-6> NA#42
43	208 DIBROMOCHLOROMETHANE <124-48-1> NA#43
44	245 1,2-DIBROMOETHANE <1060-93-4> NA#44
45	207 CHLOROBENZENE <108-90-7> NA#45
46	273 1,1,1,2-TETRACHLOROETHANE <630-20-6> NA#46

NO NAME
 47 219 ETHYLBENZENE <100-41-4> NA#47
 48 330 M,P-XYLENE <133-02-7> NA#48
 49 239 O-XYLENE <133-02-7> NA#49
 50 251 STYRENE <100-42-5> NA#50
 51 205 BROMOFORM <75-25-2> NA#51
 52 274 CIS-1,4-DICHLORO-2-BUTENE <764-71-0> NA#52
 53 275 1,2,3-TRICHLOROPROPANE <96-18-4> NA#53
 54 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> NA#54
 55 290 TRANS-1,4-DICHLORO-2-BUTENE <110-57-6> NA#55
 56 262 1,2-DIBROMO-3-CHLOROPROPANE <96-12-8> NA#56
 57 #258 D4-1,2-DICHLOROETHANE NA#57
 58 #247 BROMOFLUOROBENZENE <460-00-4> NA#58
 59 #233 DB-TOLUENE NA#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	410	5:07	1	1.000	A BB	43992.	50.000 UG/L	0.35
2	50	68	0:51	1	0.166	A BB	88667.	115.744 UG/L	0.81
3	62	77	0:58	1	0.188	A BB	114634.	123.051 UG/L	0.86
4	94	95	1:11	1	0.232	A BB	140205.	144.059 UG/L	1.01
5	64	104	1:18	1	0.254	A BB	70679.	144.564 UG/L	1.01
6	101	123	1:32	1	0.300	A BB	188225.	158.532 UG/L	1.11
7	56	181	2:16	1	0.441	A BB	78157.	2282.980 UG/L	16.02
8	96	172	2:09	1	0.420	A BB	96709.	177.519 UG/L	1.25
9	76	181	2:16	1	0.441	A BB	333585.	250.630 UG/L	1.76
10	142	182	2:16	1	0.444	A BB	365340.	231.854 UG/L	1.63
11	117	182	2:16	1	0.444	A BB	115313.	220.687 UG/L	1.55
12	85	185	2:19	1	0.451	A BB	125702.	251.987 UG/L	1.77
13	43	206	2:34	1	0.502	A BB	36738.	179.955 UG/L	1.26
14	114	542	6:46	14	1.000	A BB	162016.	50.000 UG/L	0.35
15	76	219	2:44	1	0.534	A BB	81052.	137.548 UG/L	0.97
16	84	237	2:58	1	0.578	A BB	196196.	150.783 UG/L	1.06
17	96	264	3:18	1	0.644	A BB	186701.	157.884 UG/L	1.11
18	53	280	3:30	1	0.683	A BB	390716.	1505.730 UG/L	10.57
19	63	314	3:55	1	0.766	A BB	348461.	145.228 UG/L	1.02
20	43	340	4:15	14	0.627	A BB	490189.	151.686 UG/L	1.06
21	96	384	4:48	1	0.937	A BB	226505.	146.850 UG/L	1.03
22	72	407	5:05	1	0.993	A BB	16330.	172.774 UG/L	1.21
23	83	431	5:23	1	1.051	A BB	396908.	158.089 UG/L	1.11
24	97	432	5:24	14	0.797	A BB	361703.	160.183 UG/L	1.12
25	117	448	5:36	14	0.827	A VB	333314.	153.763 UG/L	1.08
26	78	476	5:57	14	0.878	A BB	426843.	153.488 UG/L	1.08
27	62	488	6:06	1	1.190	A BB	278509.	148.748 UG/L	1.04
28	70	542	6:46	14	1.000	A BB	73351.	1522.070 UG/L	10.68
29	117	916	11:27	29	1.000	A BB	201041.	50.000 UG/L	0.35
30	130	561	7:01	14	1.035	A BB	207493.	162.956 UG/L	1.14
31	63	589	7:22	14	1.087	A BB	176263.	150.652 UG/L	1.06
32	174	604	7:33	1	1.473	A BB	215121.	179.115 UG/L	1.26
33	83	636	7:57	14	1.173	A BB	329132.	152.833 UG/L	1.07
34	63	689	8:37	14	1.271	A BB	70406.	155.513 UG/L	1.09
35	75	693	8:40	14	1.279	A BB	284481.	157.862 UG/L	1.11
36	43	729	9:07	29	0.796	M XX	141877.	162.736 UG/L	1.14
37	92	725	9:04	29	0.791	A BB	198577.	154.171 UG/L	1.08
38	75	775	9:41	14	1.430	A BB	220082.	135.742 UG/L	0.95
39	97	795	9:56	14	1.467	A BB	250560.	166.005 UG/L	1.16
40	69	803	10:02	29	0.877	A BB	350415.	146.604 UG/L	1.03
41	164	791	9:53	29	0.864	A BB	320778.	166.269 UG/L	1.17

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
42	43	848	10:36	29	0.926	A BB	136819.	157.088 UG/L	1.10
43	129	842	10:31	14	1.554	A BB	405644.	166.027 UG/L	1.17
44	107	845	10:34	14	1.559	A BB	386154.	168.239 UG/L	1.18
45	112	921	11:31	29	1.005	A BB	578171.	148.546 UG/L	1.04
46	131	943	11:47	14	1.740	A BB	347573.	170.189 UG/L	1.19
47	106	947	11:50	29	1.034	A BV	246163.	153.159 UG/L	1.07
48	106	966	12:04	29	1.055	A VB	424052.	154.852 UG/L	1.09
49	106	1021	12:46	29	1.119	A BB	354552.	154.716 UG/L	1.09
50	104	1027	12:50	29	1.121	A BB	620824.	155.950 UG/L	1.09
51	173	1048	13:06	14	1.934	A BB	364594.	171.783 UG/L	1.21
52	88	1105	13:49	14	2.039	A BB	125941.	166.499 UG/L	1.17
53	110	1138	14:13	29	1.242	A BB	128620.	155.649 UG/L	1.09
54	83	1145	14:19	29	1.250	A BB	412598.	145.547 UG/L	1.02
55	53	1152	14:24	29	1.258	A BB	99762.	145.228 UG/L	1.02
56	157	1463	18:17	29	1.597	A BB	170498.	400.953 UG/L	2.81
57	65	479	5:59	1	1.168	A BB	262031.	148.044 UG/L	1.04
58	95	1099	13:44	29	1.200	A BB	530281.	147.851 UG/L	1.04
59	98	717	8:58	29	0.783	A BB	332877.	152.578 UG/L	1.07

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:05	1.01	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	0:52	0.99	10.000	0.02	115.74	200.00	0.504	0.871	0.58
3	0:58	1.00	10.000	0.02	123.05	200.00	0.651	1.059	0.62
4	1:11	1.00	10.000	0.02	144.06	200.00	0.797	1.106	0.72
5	1:17	1.01	10.000	0.03	144.56	200.00	0.402	0.556	0.72
6	1:31	1.01	10.000	0.03	158.53	200.00	1.070	1.349	0.79
7	2:12	1.03	90.000	0.00	2282.99	2000.04	0.044	0.039	1.14
8	2:07	1.01	5.000	0.08	177.52	200.00	0.550	0.619	0.89
9	2:13	1.02	5.000	0.09	250.63	200.00	1.896	1.513	1.25
10	2:15	1.01	10.000	0.04	231.85	200.00	2.076	1.791	1.16
11	2:15	1.01	10.000	0.04	220.69	200.00	0.655	0.594	1.10
12	2:16	1.02	10.000	0.05	251.99	200.00	0.714	0.567	1.26
13	2:35	1.00	10.000	0.05	179.96	200.00	0.209	0.232	0.90
14	6:43	1.01	5.000	0.20	50.00	50.00	1.000	1.000	1.00
15	2:43	1.01	10.000	0.05	137.55	200.00	0.461	0.670	0.69
16	2:56	1.01	5.000	0.12	150.78	200.00	1.115	1.479	0.75
17	3:16	1.01	5.000	0.13	157.88	200.00	1.061	1.344	0.79
18	3:29	1.00	120.000	0.01	1505.73	2000.04	0.222	0.299	0.75
19	3:54	1.01	5.000	0.15	145.23	200.00	1.980	2.727	0.73
20	4:13	1.01	10.000	0.06	151.69	200.00	0.756	0.997	0.76
21	4:46	1.01	5.000	0.19	146.85	200.00	1.287	1.753	0.73
22	5:04	1.00	10.000	0.10	172.77	200.00	0.093	0.107	0.86
23	5:20	1.01	5.000	0.21	158.09	200.00	2.256	2.854	0.79
24	5:21	1.01	5.000	0.16	160.18	200.00	0.558	0.697	0.80
25	5:33	1.01	5.000	0.17	153.76	200.00	0.514	0.669	0.77
26	5:54	1.01	5.000	0.18	153.49	200.00	0.659	0.858	0.77
27	6:04	1.01	5.000	0.24	148.75	200.00	1.583	2.128	0.74
28	6:46	1.00	100.000	0.01	1522.07	2000.04	0.011	0.015	0.76
29	11:23	1.01	5.000	0.20	50.00	50.00	1.000	1.000	1.00
30	6:58	1.01	5.000	0.21	162.96	200.00	0.320	0.393	0.81
31	7:19	1.01	5.000	0.22	150.65	200.00	0.272	0.361	0.75
32	7:30	1.01	5.000	0.29	179.11	200.00	1.223	1.365	0.90
33	7:53	1.01	5.000	0.23	152.83	200.00	0.508	0.665	0.76
34	8:33	1.01	10.000	0.13	155.51	200.00	0.109	0.140	0.78
35	8:36	1.01	5.000	0.26	157.86	200.00	0.439	0.556	0.79
36	9:05	1.00	15.000	0.05	162.74	200.00	0.176	0.217	0.81

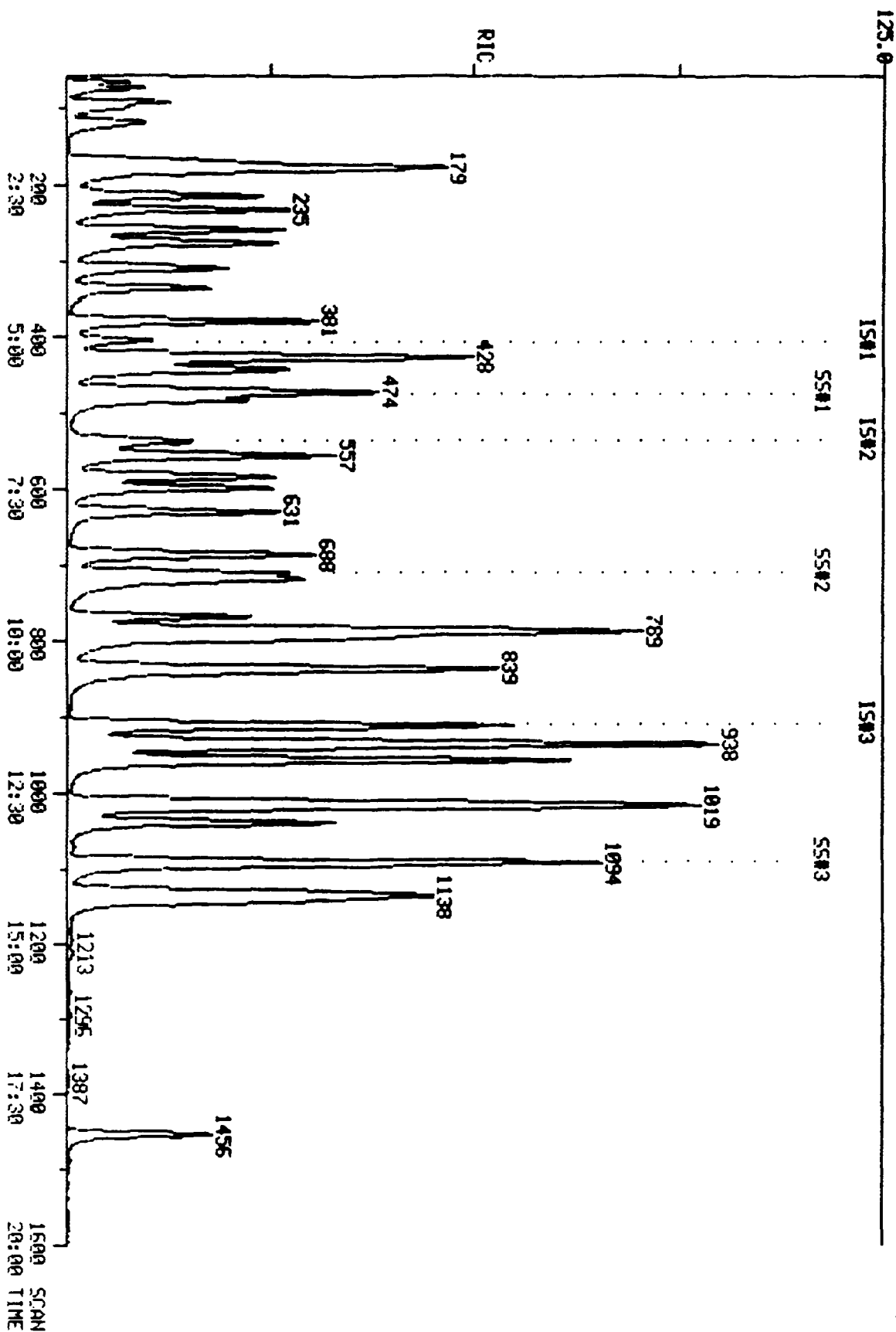
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
37	9:00	1.01	5.000	0.16	154.17	200.00	0.247	0.320	0.77
38	9:37	1.01	5.000	0.29	135.74	200.00	0.340	0.500	0.68
39	9:53	1.01	5.000	0.29	166.00	200.00	0.387	0.466	0.83
40	9:59	1.01	10.000	0.09	146.60	200.00	0.436	0.594	0.73
41	9:50	1.01	5.000	0.17	166.27	200.00	0.399	0.480	0.83
42	10:34	1.00	15.000	0.06	157.09	200.00	0.170	0.217	0.79
43	10:28	1.01	5.000	0.31	166.03	200.00	0.626	0.754	0.83
44	10:31	1.00	5.000	0.31	168.24	200.00	0.596	0.708	0.84
45	11:25	1.01	5.000	0.20	148.55	200.00	0.719	0.968	0.74
46	11:41	1.01	5.000	0.35	170.19	200.00	0.536	0.630	0.85
47	11:46	1.01	5.000	0.21	153.16	200.00	0.306	0.400	0.77
48	11:59	1.01	5.000	0.21	154.85	200.00	0.527	0.681	0.77
49	12:40	1.01	5.000	0.22	154.72	200.00	0.441	0.570	0.77
50	12:46	1.01	5.000	0.22	155.95	200.00	0.772	0.990	0.78
51	13:01	1.01	5.000	0.39	171.78	200.00	0.563	0.655	0.86
52	13:44	1.01	15.000	0.14	166.50	200.00	0.194	0.233	0.83
53	14:09	1.01	15.000	0.08	155.65	200.00	0.160	0.206	0.78
54	14:13	1.01	5.000	0.25	145.55	200.00	0.513	0.705	0.73
55	14:19	1.01	15.000	0.08	145.23	200.00	0.124	0.171	0.73
56	18:12	1.00	10.000	0.16	400.95	400.00	0.106	0.106	1.00
57	5:56	1.01	5.000	0.23	148.04	200.00	1.489	2.012	0.74
58	13:40	1.01	5.000	0.24	147.85	200.00	0.659	0.892	0.74
59	8:54	1.01	5.000	0.16	152.58	200.00	0.414	0.543	0.76

RIC
12/28/89 10:34:00
SAMPLE: SML UST0200 #1915 ON#13
CONDOS.:

COMPUCHEM LABS

COMPUCHEM DATA: C5891228A13 SCANS 59 TO 1600

603520.



QUANTITATION REPORT FILE: CS891228A13
DATA: CS891228A13.TI
12/28/89 10:34:00
SAMPLE: 5ML VSTD200 #1915 DN#13
CONDS. :
SUBMITTED BY: 13 ANALYST: 1492

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> NA#1
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3	231 VINYL CHLORIDE <75-01-4> NA#3
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5	209 CHLOROETHANE <75-00-3> NA#5
6	230 TRICHLOROFLUOROMETHANE <75-69-4> NA#6
7	201 ACROLEIN <107-02-8> NA#7
8	216 1,1-DICHLOROETHENE <75-35-4> NA#8
9	254 CARBON DISULFIDE <75-15-0> NA#9
10	285 IODOMETHANE <74-88-4> NA#10
11	297 1,1,1-TRICHLORO-2,2,2-TRIFLUOROETHANE <354-58-5> NA#11
12	266 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE <76-13-1> NA#12
13	252 ACETONE (2-PROPANONE) <67-64-1> NA#13
14	*248 1,4-DIFLUOROBENZENE (IS) <540-36-3> NA#14
15	298 3-CHLOROPROPENE <107-05-1> NA#15
16	222 METHYLENE CHLORIDE <75-09-2> NA#16
17	226 TRANS-1,2-DICHLOROETHENE <156-60-5> NA#17
18	202 ACRYLONITRILE <107-13-1> NA#18
19	214 1,1-DICHLOROETHANE <75-34-3> NA#19
20	257 VINYL ACETATE <108-05-4> NA#20
21	237 CIS-1,2-DICHLOROETHENE <156-59-2> NA#21
22	253 2-BUTANONE <78-93-3> NA#22
23	211 CHLOROFORM <67-66-2> NA#23
24	227 1,1,1-TRICHLOROETHANE <71-55-6> NA#24
25	206 CARBON TETRACHLORIDE <56-23-5> NA#25
26	203 BENZENE <71-43-2> NA#26
27	215 1,2-DICHLOROETHANE <107-06-2> NA#27
28	272 CROTONALDEHYDE <4170-30-3> NA#28
29	*270 D5-CHLOROBENZENE (IS) <XX-XX-X> NA#29
30	229 TRICHLOROETHENE <79-01-6> NA#30
31	217 1,2-DICHLOROPROPANE <78-87-5> NA#31
32	286 DIBROMOMETHANE <74-95-3> NA#32
33	212 BROMODICHLOROMETHANE <75-27-4> NA#33
34	210 2-CHLOROETHYL VINYL ETHER <110-75-8> NA#34
35	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> NA#35
36	256 4-METHYL-2-PENTANONE <108-01-1> NA#36
37	225 TOLUENE <108-88-3> NA#37
38	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> NA#38
39	228 1,1,2-TRICHLOROETHANE <79-00-5> NA#39
40	287 ETHYLMETHACRYLATE <96-18-4> NA#40
41	224 TETRACHLOROETHENE <127-18-4> NA#41
42	255 2-HEXANONE <591-78-6> NA#42
43	208 DIBROMOCHLOROMETHANE <124-48-1> NA#43
44	245 1,2-DIBROMOETHANE <1060-93-4> NA#44
45	207 CHLOROBENZENE <108-90-7> NA#45
46	273 1,1,1,2-TETRACHLOROETHANE <630-20-6> NA#46

NO NAME
 47 219 ETHYLBENZENE <100-41-4> NA#47
 48 330 M,P-XYLENE <133-02-7> NA#48
 49 239 O-XYLENE <133-02-7> NA#49
 50 251 STYRENE <100-42-5> NA#50
 51 205 BROMOFORM <75-25-2> NA#51
 52 274 CIS-1,4-DICHLORO-2-BUTENE <764-71-0> NA#52
 53 275 1,2,3-TRICHLOROPROPANE <96-18-4> NA#53
 54 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> NA#54
 55 290 TRANS-1,4-DICHLORO-2-BUTENE <110-57-6> NA#55
 56 262 1,2-DIBROMO-3-CHLOROPROPANE <96-12-8> NA#56
 57 #258 D4-1,2-DICHLOROETHANE NA#57
 58 #247 BROMOFLUOROBENZENE <460-00-4> NA#58
 59 #233 DB-TOLUENE NA#59

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	UG/L	%TOT
1	128	407	5:05	1	1.000	A BB	45838.	50.000	UG/L	0.29
2	50	69	0:52	1	0.170	A BB	159642.	200.000	UG/L	1.18
3	62	77	0:58	1	0.189	A BB	194140.	200.000	UG/L	1.18
4	94	95	1:11	1	0.233	A BB	202817.	200.000	UG/L	1.18
5	64	103	1:17	1	0.253	A BB	101885.	200.000	UG/L	1.18
6	101	122	1:31	1	0.300	A BB	247425.	200.000	UG/L	1.18
7	56	176	2:12	1	0.432	A BB	71343.	2000.030	UG/L	11.80
8	96	170	2:07	1	0.418	A BB	113528.	200.000	UG/L	1.18
9	76	178	2:13	1	0.437	A BB	277367.	200.000	UG/L	1.18
10	142	180	2:15	1	0.442	A BB	328371.	200.000	UG/L	1.18
11	117	180	2:15	1	0.442	A BB	108889.	200.000	UG/L	1.18
12	85	182	2:16	1	0.447	A BB	103955.	200.000	UG/L	1.18
13	43	207	2:35	1	0.509	M XX	42544.	200.000	UG/L	1.18
14	114	538	6:43	14	1.000	A BB	167360.	50.000	UG/L	0.29
15	76	217	2:43	1	0.533	A BB	122798.	200.000	UG/L	1.18
16	84	235	2:56	1	0.577	A BB	271157.	200.000	UG/L	1.18
17	96	262	3:16	1	0.644	A BB	246428.	200.000	UG/L	1.18
18	53	279	3:29	1	0.686	A BB	540759.	2000.030	UG/L	11.80
19	63	312	3:54	1	0.767	A BB	500017.	200.000	UG/L	1.18
20	43	338	4:13	14	0.628	A BB	668435.	200.000	UG/L	1.18
21	96	381	4:46	1	0.936	A BB	321429.	200.000	UG/L	1.18
22	72	405	5:04	1	0.995	A BB	19696.	200.000	UG/L	1.18
23	83	427	5:20	1	1.049	A BB	523202.	200.000	UG/L	1.18
24	97	428	5:21	14	0.796	A BB	467066.	200.000	UG/L	1.18
25	117	444	5:33	14	0.825	A VB	448377.	200.000	UG/L	1.18
26	78	472	5:54	14	0.877	A BB	575221.	200.000	UG/L	1.18
27	62	485	6:04	1	1.192	A BB	390184.	200.000	UG/L	1.18
28	70	541	6:46	14	1.006	A BB	99683.	2000.030	UG/L	11.80
29	117	911	11:23	29	1.000	A BB	193949.	50.000	UG/L	0.29
30	130	557	6:58	14	1.035	A BB	263376.	200.000	UG/L	1.18
31	63	585	7:19	14	1.087	A BB	242008.	200.000	UG/L	1.18
32	174	600	7:30	1	1.474	A BB	250284.	200.000	UG/L	1.18
33	83	631	7:53	14	1.173	A BB	445447.	200.000	UG/L	1.18
34	63	684	8:33	14	1.271	A BB	93645.	200.000	UG/L	1.18
35	75	688	8:36	14	1.279	A BB	372749.	200.000	UG/L	1.18
36	43	727	9:05	29	0.798	A BB	168214.	200.000	UG/L	1.18
37	92	720	9:00	29	0.790	A BB	248520.	200.000	UG/L	1.18
38	75	770	9:37	14	1.431	A BB	335360.	200.000	UG/L	1.18
39	97	791	9:53	14	1.470	A BB	312200.	200.000	UG/L	1.18
40	69	799	9:59	29	0.877	A BB	461178.	200.000	UG/L	1.18
41	164	787	9:50	29	0.864	A BB	372243.	200.000	UG/L	1.18

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
42	43	845	10:34	29	0.928	A BB	168049.	200.000 UG/L	1.18
43	129	837	10:28	14	1.556	A BB	505368.	200.000 UG/L	1.18
44	107	841	10:31	14	1.563	A BB	474763.	200.000 UG/L	1.18
45	112	914	11:25	29	1.003	A BB	750980.	200.000 UG/L	1.18
46	131	935	11:41	14	1.738	A BB	422433.	200.000 UG/L	1.18
47	106	941	11:46	29	1.033	A BV	310108.	200.000 UG/L	1.18
48	106	959	11:59	29	1.053	A VB	528368.	200.000 UG/L	1.18
49	106	1014	12:40	29	1.113	A BB	442157.	200.000 UG/L	1.18
50	104	1021	12:46	29	1.121	A BB	768099.	200.000 UG/L	1.18
51	173	1042	13:01	14	1.937	A BB	439008.	200.000 UG/L	1.18
52	88	1099	13:44	14	2.043	A BB	156458.	200.000 UG/L	1.18
53	110	1132	14:09	29	1.243	A BB	159439.	200.000 UG/L	1.18
54	83	1138	14:13	29	1.249	A BB	546961.	200.000 UG/L	1.18
55	53	1146	14:19	29	1.258	A BB	132540.	200.000 UG/L	1.18
56	157	1456	18:12	29	1.598	A BB	164092.	399.999 UG/L	2.36
57	65	475	5:56	1	1.167	A BB	368845.	200.000 UG/L	1.18
58	95	1093	13:40	29	1.200	A BB	692016.	200.000 UG/L	1.18
59	98	712	8:54	29	0.782	A BB	420945.	200.000 UG/L	1.18

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	5:05	1.00	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	0:52	1.00	10.000	0.02	200.00	200.00	0.871	0.871	1.00
3	0:58	1.00	10.000	0.02	200.00	200.00	1.059	1.059	1.00
4	1:11	1.00	10.000	0.02	200.00	200.00	1.106	1.106	1.00
5	1:17	1.00	10.000	0.03	200.00	200.00	0.556	0.556	1.00
6	1:31	1.00	10.000	0.03	200.00	200.00	1.349	1.349	1.00
7	2:12	1.00	90.000	0.00	2000.04	2000.04	0.039	0.039	1.00
8	2:07	1.00	5.000	0.08	200.00	200.00	0.619	0.619	1.00
9	2:13	1.00	5.000	0.09	200.00	200.00	1.513	1.513	1.00
10	2:15	1.00	10.000	0.04	200.00	200.00	1.791	1.791	1.00
11	2:15	1.00	10.000	0.04	200.00	200.00	0.594	0.594	1.00
12	2:16	1.00	10.000	0.04	200.00	200.00	0.567	0.567	1.00
13	2:35	1.00	10.000	0.05	200.00	200.00	0.232	0.232	1.00
14	6:43	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
15	2:43	1.00	10.000	0.05	200.00	200.00	0.670	0.670	1.00
16	2:56	1.00	5.000	0.12	200.00	200.00	1.479	1.479	1.00
17	3:16	1.00	5.000	0.13	200.00	200.00	1.344	1.344	1.00
18	3:29	1.00	120.000	0.01	2000.04	2000.04	0.295	0.295	1.00
19	3:54	1.00	5.000	0.15	200.00	200.00	2.727	2.727	1.00
20	4:13	1.00	10.000	0.06	200.00	200.00	0.997	0.997	1.00
21	4:46	1.00	5.000	0.19	200.00	200.00	1.753	1.753	1.00
22	5:04	1.00	10.000	0.10	200.00	200.00	0.107	0.107	1.00
23	5:20	1.00	5.000	0.21	200.00	200.00	2.854	2.854	1.00
24	5:21	1.00	5.000	0.16	200.00	200.00	0.697	0.697	1.00
25	5:33	1.00	5.000	0.17	200.00	200.00	0.669	0.669	1.00
26	5:54	1.00	5.000	0.18	200.00	200.00	0.858	0.858	1.00
27	6:04	1.00	5.000	0.24	200.00	200.00	2.128	2.128	1.00
28	6:46	1.00	100.000	0.01	2000.04	2000.04	0.015	0.015	1.00
29	11:23	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
30	6:58	1.00	5.000	0.21	200.00	200.00	0.393	0.393	1.00
31	7:19	1.00	5.000	0.22	200.00	200.00	0.361	0.361	1.00
32	7:30	1.00	5.000	0.29	200.00	200.00	1.365	1.365	1.00
33	7:53	1.00	5.000	0.23	200.00	200.00	0.665	0.665	1.00
34	8:33	1.00	10.000	0.13	200.00	200.00	0.140	0.140	1.00
35	8:36	1.00	5.000	0.26	200.00	200.00	0.556	0.556	1.00
36	9:05	1.00	15.000	0.05	200.00	200.00	0.217	0.217	1.00

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
37	9:00	1.00	5.000	0.16	200.00	200.00	0.320	0.320	1.00
38	9:37	1.00	5.000	0.29	200.00	200.00	0.500	0.500	1.00
39	9:53	1.00	5.000	0.29	200.00	200.00	0.466	0.466	1.00
40	9:59	1.00	10.000	0.09	200.00	200.00	0.594	0.594	1.00
41	9:50	1.00	5.000	0.17	200.00	200.00	0.480	0.480	1.00
42	10:34	1.00	15.000	0.06	200.00	200.00	0.217	0.217	1.00
43	10:28	1.00	5.000	0.31	200.00	200.00	0.754	0.754	1.00
44	10:31	1.00	5.000	0.31	200.00	200.00	0.708	0.708	1.00
45	11:25	1.00	5.000	0.20	200.00	200.00	0.968	0.968	1.00
46	11:41	1.00	5.000	0.35	200.00	200.00	0.630	0.630	1.00
47	11:46	1.00	5.000	0.21	200.00	200.00	0.400	0.400	1.00
48	11:59	1.00	5.000	0.21	200.00	200.00	0.681	0.681	1.00
49	12:40	1.00	5.000	0.22	200.00	200.00	0.570	0.570	1.00
50	12:46	1.00	5.000	0.22	200.00	200.00	0.990	0.990	1.00
51	13:01	1.00	5.000	0.39	200.00	200.00	0.655	0.655	1.00
52	13:44	1.00	15.000	0.14	200.00	200.00	0.233	0.233	1.00
53	14:09	1.00	15.000	0.08	200.00	200.00	0.206	0.206	1.00
54	14:13	1.00	5.000	0.25	200.00	200.00	0.705	0.705	1.00
55	14:19	1.00	15.000	0.08	200.00	200.00	0.171	0.171	1.00
56	18:12	1.00	10.000	0.16	400.00	400.00	0.106	0.106	1.00
57	5:56	1.00	5.000	0.23	200.00	200.00	2.012	2.012	1.00
58	13:40	1.00	5.000	0.24	200.00	200.00	0.892	0.892	1.00
59	8:54	1.00	5.000	0.16	200.00	200.00	0.543	0.543	1.00

COMPUCHEM LABORATORIES, INC.
GC/MS ANALYSIS LOG

INITIAL TIME OF TUNE 10:18 SHIFT(S) (A) X (B) _____ (C) _____
TIME TUNE EXPIRES 22:18 DATE 12/28/89
ANALYSIS TYPE NAST

RUN LOG

PREVENTIVE MAINTENANCE _____

FILE NAME	DATE	TIME	EPA ID	CASE NO	STD ID #	AMOUNT INJECTED	CHEMIST	COMMENTS (Lot #, s, Disposition, Etc.)
BE891228A13	11/11	10:18	BFB			2ul	1492	H7008
CS891228A13	11/11	10:34	VSTD0200		1915	5ul	1492	
CT891228A13	11/11	11:08	VSTD0200		1911	5ul	1492	
CU891228A13	11/11	11:52	VSTD150		1914	5ul	1492	
CV891228A13	11/11	12:30	VSTD050		1912	5ul	1492	
CW891228A13	11/11	13:04	VSTD100		1913	5ul	1492	
			NAST MP 6000					

VERIFIED WJL/STH
SUPERVISOR APPROVAL 1452 *[Signature]*

6A
VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: COMPUCHEM LABS Contract: (2-88)-REVS

Lab Code: COMPU Case No.: 18756 SAS No.: _____ SDG No.: 07

Instrument ID: 19 Calibration Date(s): 12/12/89 12/13/89

Matrix: (soil/water) SOIL Level: (low/med) LOW Column: (pack/cap) CAP

Min \overline{RRF} for SPCC(#) = 0.300 (0.250 for Bromoform) Max %RSD for CCC(*) = 30.0%

LAB FILE ID: RRF20 = GS891213C19 RRF50 = GT891212B19
 RRF100 = GW891212B19 RRF150 = GU891213C19 RRF200 = GU891212B19

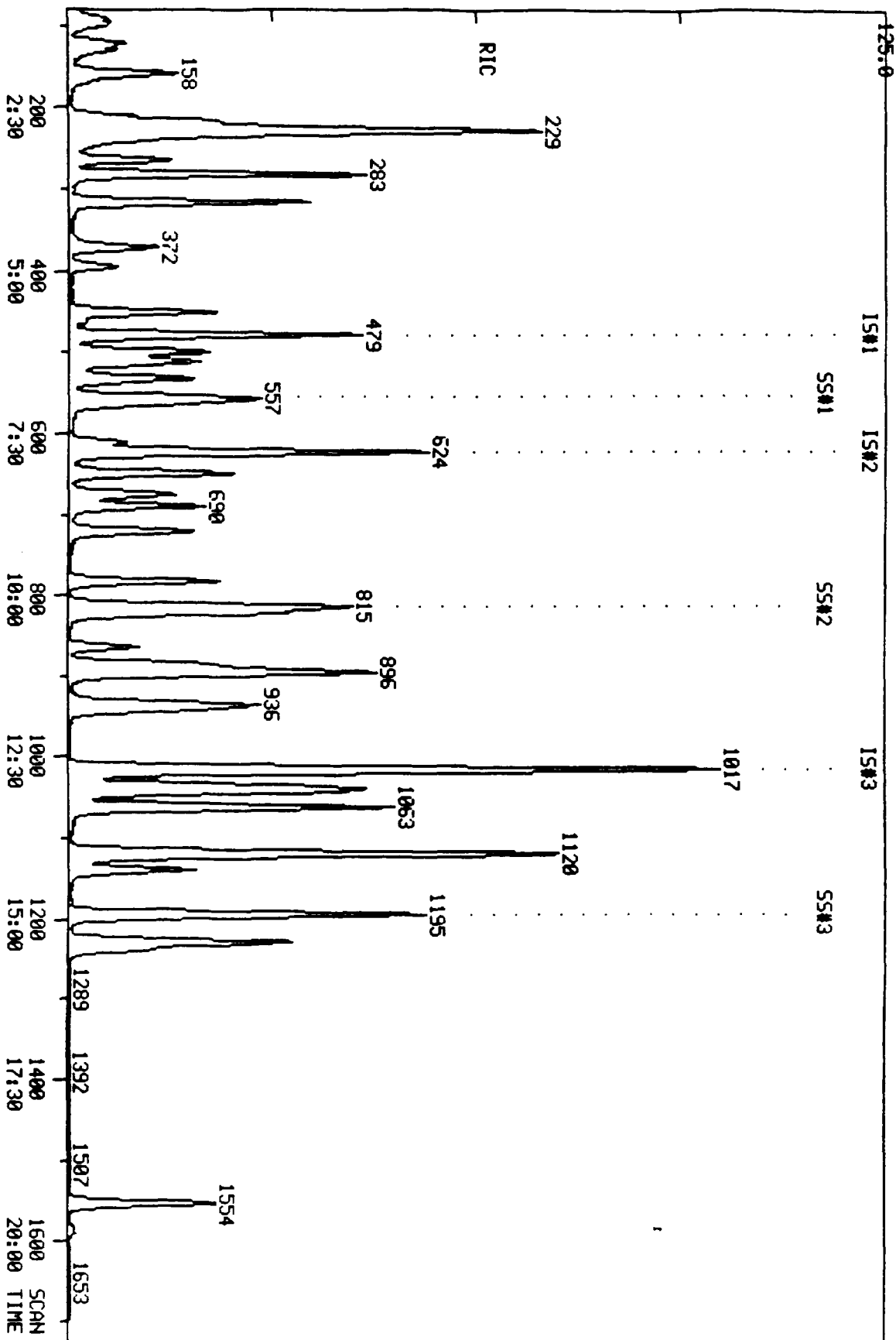
COMPOUND	RRF20	RRF50	RRF100	RRF150	RRF200	\overline{RRF}	% RSD
Chloromethane	1.440	1.182	1.251	1.328	1.407	1.322	8.1#
Bromomethane	1.727	1.457	1.288	1.417	1.415	1.461	11.1
Vinyl Chloride	1.973	1.406	1.526	1.614	1.586	1.621	13.1*
Chloroethane	0.889	0.850	0.641	0.645	0.705	0.746	15.6
Methylene Chloride	4.228	2.065	1.924	1.732	1.845	2.359	44.6
Acetone	0.659	0.385	0.357	0.359	0.276	0.407	36.0
Carbon Disulfide	4.886	5.046	4.423	3.515	3.811	4.336	15.3
1,1-Dichloroethene	1.604	1.458	1.509	1.290	1.504	1.473	7.8*
1,1-Dichloroethane	3.309	3.038	3.372	3.207	3.488	3.283	5.2#
1,2-Dichloroethene (total)	3.715	3.247	3.438	3.277	3.466	3.429	5.4
Chloroform	3.900	3.567	3.772	3.644	3.800	3.737	3.5*
1,2-Dichloroethane	2.844	2.476	2.640	2.683	2.590	2.647	5.1
2-Butanone	0.147	0.111	0.151	0.163	0.127	0.140	14.8
1,1,1-Trichloroethane	0.863	0.807	0.883	0.900	0.834	0.857	4.4
Carbon Tetrachloride	0.729	0.743	0.813	0.826	0.796	0.781	5.5
Vinyl Acetate	0.654	0.637	0.745	0.819	0.645	0.700	11.4
Bromodichloromethane	0.795	0.792	0.930	1.018	0.858	0.879	10.9
1,2-Dichloropropane	0.384	0.330	0.361	0.387	0.330	0.358	7.8*
cis-1,3-Dichloropropene	0.884	0.793	0.968	1.022	0.873	0.908	9.8
Trichloroethene	0.500	0.413	0.465	0.472	0.436	0.457	7.4
Dibromochloromethane	0.695	0.647	0.737	0.806	0.672	0.711	8.8
1,1,2-Trichloroethane	0.422	0.346	0.422	0.466	0.371	0.405	11.7
Benzene	0.946	0.821	0.907	0.894	0.800	0.874	7.0
Trans-1,3-Dichloropropene	0.429	0.386	0.476	0.532	0.461	0.457	11.9
Bromoform	0.610	0.564	0.683	0.779	0.623	0.652	12.7#
4-Methyl-2-Pentanone	0.408	0.336	0.451	0.503	0.388	0.417	15.2
2-Hexanone	0.295	0.225	0.385	0.407	0.289	0.320	23.4
Tetrachloroethene	0.592	0.514	0.565	0.543	0.527	0.548	5.7
1,1,2,2-Tetrachloroethane	0.549	0.553	0.691	0.767	0.650	0.642	14.5#
Toluene	0.737	0.657	0.760	0.746	0.724	0.725	5.5*
Chlorobenzene	1.064	0.921	1.053	1.060	1.011	1.022	5.9#
Ethylbenzene	0.482	0.436	0.491	0.471	0.460	0.468	4.6*
Styrene	1.290	1.194	1.212	1.166	1.071	1.187	6.7
Total Xylenes	1.606	1.573	1.563	1.490	1.405	1.527	5.3
=====							
Toluene-d8	1.270	1.119	1.291	1.129	1.101	1.182	7.7
BFB	1.061	0.991	1.103	0.983	0.937	1.015	6.5
1,2-Dichloroethane-d4	2.700	2.601	2.731	2.493	2.451	2.595	4.7

RIC
 12/13/89 0:31:00
 SAMPLE: 10ML EPA STANDARD UST0020 (1906) DN#19
 COND5.:

COMPUchem LABS

COMPUchem DATA: GS891213C19 SCANS 80 TO 1700
 OUT OF 80 TO 1700

104160.



QUANTITATION REPORT FILE: QS891213C19
DATA: QS891213C19.TI ✓
12/13/89 0:31:00 ✓
SAMPLE: 10ML EPA STANDARD VSTD020 (1906) DN#19
CONDS.:
SUBMITTED BY: 19 ANALYST: 1422 ✓

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
RESP. FAC. FROM LIBRARY ENTRY ✓

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1
2	221 CHLOROMETHANE <74-87-3> RO#2
3	231 VINYL CHLORIDE <75-01-4> RO#3
4	220 BROMOMETHANE <78-83-9> RO#4
5	209 CHLOROETHANE <75-00-3> RO#5
6	230 TRICHLOROFLUOROMETHANE <75-69-4> RO#6
7	201 ACROLEIN <107-02-8> RO#7
8	216 1,1-DICHLOROETHENE <75-35-4> RO#8
9	254 CARBON DISULFIDE <75-15-0> RO#9
10	285 IODOMETHANE <74-88-4> RO#10
11	297 1,1,1-TRICHLORO-2,2,2-TRIFLUOROETHANE <354-58-5> RO#11
12	266 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE <76-13-1> RO#12
13	252 ACETONE (2-PROPANONE) <67-64-1> RO#13
14	*248 1,4-DIFLUOROBENZENE (IS) <340-36-3> RO#14
15	298 3-CHLOROPROPENE <107-05-1> RO#15
16	222 METHYLENE CHLORIDE <75-09-2> RO#16
17	226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17
18	202 ACRYLONITRILE <107-13-1> RO#18
19	214 1,1-DICHLOROETHANE <75-34-3> RO#19
20	257 VINYL ACETATE <108-05-4> RO#20
21	237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21
22	253 2-BUTANONE <78-93-3> RO#22
23	211 CHLOROFORM <67-66-2> RO#23
24	227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24
25	206 CARBON TETRACHLORIDE <56-23-5> RO#25
26	203 BENZENE <71-43-2> RO#26
27	215 1,2-DICHLOROETHANE <107-06-2> RO#27
28	272 CROTONALDEHYDE <4170-30-3> RO#28
29	*270 D5-CHLOROBENZENE (IS) RO#29
30	229 TRICHLOROETHENE <79-01-6> RO#30
31	217 1,2-DICHLOROPROPANE <78-87-5> RO#31
32	286 DIBROMOMETHANE <74-95-3> RO#32
33	212 BROMODICHLOROMETHANE <75-27-4> RO#33
34	210 2-CHLOROETHYL VINYL ETHER <110-75-8> RO#34
35	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35
36	256 4-METHYL-2-PENTANONE <108-01-1> RO#36
37	225 TOLUENE <108-88-3> RO#37
38	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38
39	228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39
40	287 ETHYLMETHACRYLATE <96-18-4> RO#40
41	224 TETRACHLOROETHENE <127-18-4> RO#41
42	255 2-HEXANONE <591-78-6> RO#42
43	208 DIBROMOCHLOROMETHANE <124-48-1> RO#43
44	245 1,2-DIBROMOMETHANE <1060-93-4> RO#44
45	207 CHLOROBENZENE <108-90-7> RO#45
46	273 1,1,1,2-TETRACHLOROETHANE <630-20-6> RO#46

NO NAME
 47 219 ETHYLBENZENE <100-41-4> RO#47
 48 330 M,P-XYLENE <133-02-7> RO#48
 49 239 O-XYLENE <133-02-7> RO#49
 50 251 STYRENE <100-42-5> RO#50
 51 205 BROMOFORM <75-25-2> RO#51
 52 274 CIS-1,4-DICHLORO-2-BUTENE <764-71-0> RO#52
 53 275 1,2,3-TRICHLOROPROPANE <96-18-4> RO#53
 54 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54
 55 290 TRANS-1,4-DICHLORO-2-BUTENE <110-57-6> RO#55
 56 262 1,2-DIBROMO-3-CHLOROPROPANE <96-12-8> RO#56
 57 #258 D4-1,2-DICHLOROETHANE RO#57
 58 #247 BROMOFLUOROBENZENE <460-00-4> RO#58
 59 #233 D8-TOLUENE RO#59

66 12-13-89
 UES
 12-13-89

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	479	5:59	1	1.000	A BB	36628.	50.000 UG/KG	2.23
2	50	90	1:07	1	0.188	A BB	21095.	24.364 UG/KG	1.08
3	62	101	1:16	1	0.211	A BB	28901.	28.057 UG/KG	1.25
4	94	122	1:31	1	0.255	A BV	25299.	23.702 UG/KG	1.05
5	64	130	1:37	1	0.271	A BB	13028.	20.929 UG/KG	0.93
6	101	160	2:00	1	0.334	A BB	57856.	19.617 UG/KG	0.87
7	56	210	2:37	1	0.438	A BV	18802.	263.679 UG/KG	11.73
8	96	217	2:43	1	0.453	A BB	23503.	22.000 UG/KG	0.98
9	76	230	2:52	1	0.480	A BB	71591.	19.366 UG/KG	0.86
10	142	227	2:50	1	0.474	A BB	80851.	22.128 UG/KG	0.98
11	117	227	2:50	1	0.474	A BB	23447.	22.511 UG/KG	1.00
12	85	232	2:54	1	0.484	A BB	26377.	22.484 UG/KG	1.00
13	43	233	2:55	1	0.486	A BB	9660.	34.223 UG/KG	1.52
14	114	624	7:48	14	1.000	A BB	146288.	50.000 UG/KG	2.23
15	76	265	3:19	1	0.553	A BB	12206.	21.454 UG/KG	0.95
16	84	283	3:32	1	0.591	A BB	61944.	40.946 UG/KG	1.82
17	96	317	3:58	1	0.662	A BV	23667.	21.827 UG/KG	0.97
18	53	317	3:58	1	0.662	A BB	38263.	245.603 UG/KG	10.93
19	63	372	4:39	1	0.777	A BB	48483.	21.783 UG/KG	0.97
20	43	396	4:57	14	0.635	A BB	38271.	20.530 UG/KG	0.91
21	96	450	5:37	1	0.939	A BB	30761.	23.766 UG/KG	1.06
22	72	464	5:48	1	0.969	A BB	2153.	26.381 UG/KG	1.17
23	83	501	6:16	1	1.046	A BB	57140.	21.868 UG/KG	0.97
24	97	512	6:24	14	0.821	A BB	50512.	21.399 UG/KG	0.95
25	117	533	6:40	14	0.854	A BB	42668.	19.638 UG/KG	0.87
26	78	558	6:58	14	0.894	A BB	55377.	23.059 UG/KG	1.03
27	62	565	7:04	1	1.180	A BB	41667.	22.972 UG/KG	1.02
28	70	611	7:38	14	0.979	A BV	12535.	379.487 UG/KG	16.89
29	117	1016	12:42	29	1.000	A BB	167999.	50.000 UG/KG	2.23
30	130	650	8:07	14	1.042	A BB	29278.	24.240 UG/KG	1.08
31	63	677	8:28	14	1.085	A BB	22445.	23.273 UG/KG	1.04
32	174	690	8:37	1	1.441	A BB	29400.	21.777 UG/KG	0.97
33	83	722	9:01	14	1.157	A BB	46543.	20.088 UG/KG	0.89
34	63	NOT FOUND							
35	75	783	9:47	14	1.255	A BB	51724.	22.306 UG/KG	0.99
36	43	815	10:11	29	0.802	A BB	27447.	24.316 UG/KG	1.08
37	92	823	10:17	29	0.810	A BB	49493.	22.431 UG/KG	1.00
38	75	864	10:48	14	1.385	A BB	25090.	22.241 UG/KG	0.99
39	97	885	11:04	14	1.418	A BB	24691.	24.369 UG/KG	1.08
40	69	893	11:10	29	0.879	A BB	42102.	22.777 UG/KG	1.01
41	164	897	11:13	29	0.883	A BB	39807.	23.030 UG/KG	1.02

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
42	43	933	11:40	29	0.918	A BB	19816.	26.231 UG/KG	1.17
43	129	936	11:42	14	1.500	A BB	40694.	21.508 UG/KG	0.96
44	107	943	11:47	14	1.511	A BB	42115.	21.445 UG/KG	0.95
45	112	1019	12:44	29	1.003	A BB	71476.	23.086 UG/KG	1.03
46	131	1037	12:58	14	1.662	A BB	39928.	20.605 UG/KG	0.92
47	106	1042	13:01	29	1.026	A BV	32367.	22.097 UG/KG	0.98
48	106	1062	13:16	29	1.045	A VB	58855.	20.435 UG/KG	0.91
49	106	1118	13:58	29	1.100	A BB	49068.	20.401 UG/KG	0.91
50	104	1122	14:01	29	1.104	A BB	86661.	21.608 UG/KG	0.96
51	173	1141	14:16	14	1.829	A BB	35672.	21.636 UG/KG	0.96
52	88	1193	14:55	14	1.912	A BB	12780.	20.112 UG/KG	0.89
53	110	1227	15:20	29	1.208	A BB	13668.	21.669 UG/KG	0.96
54	83	1231	15:23	29	1.212	A BB	36908.	19.880 UG/KG	0.88
55	53	1239	15:29	29	1.219	A BB	10760.	19.543 UG/KG	0.87
56	157	1554	19:25	29	1.530	A BB	28462.	48.047 UG/KG	2.14
57	65	555	6:56	1	1.159	A BB	39560.	20.759 UG/KG	0.92
58	95	1195	14:56	29	1.176	A BB	71276.	21.408 UG/KG	0.95
59	98	815	10:11	29	0.802	A BB	85365.	22.714 UG/KG	1.01

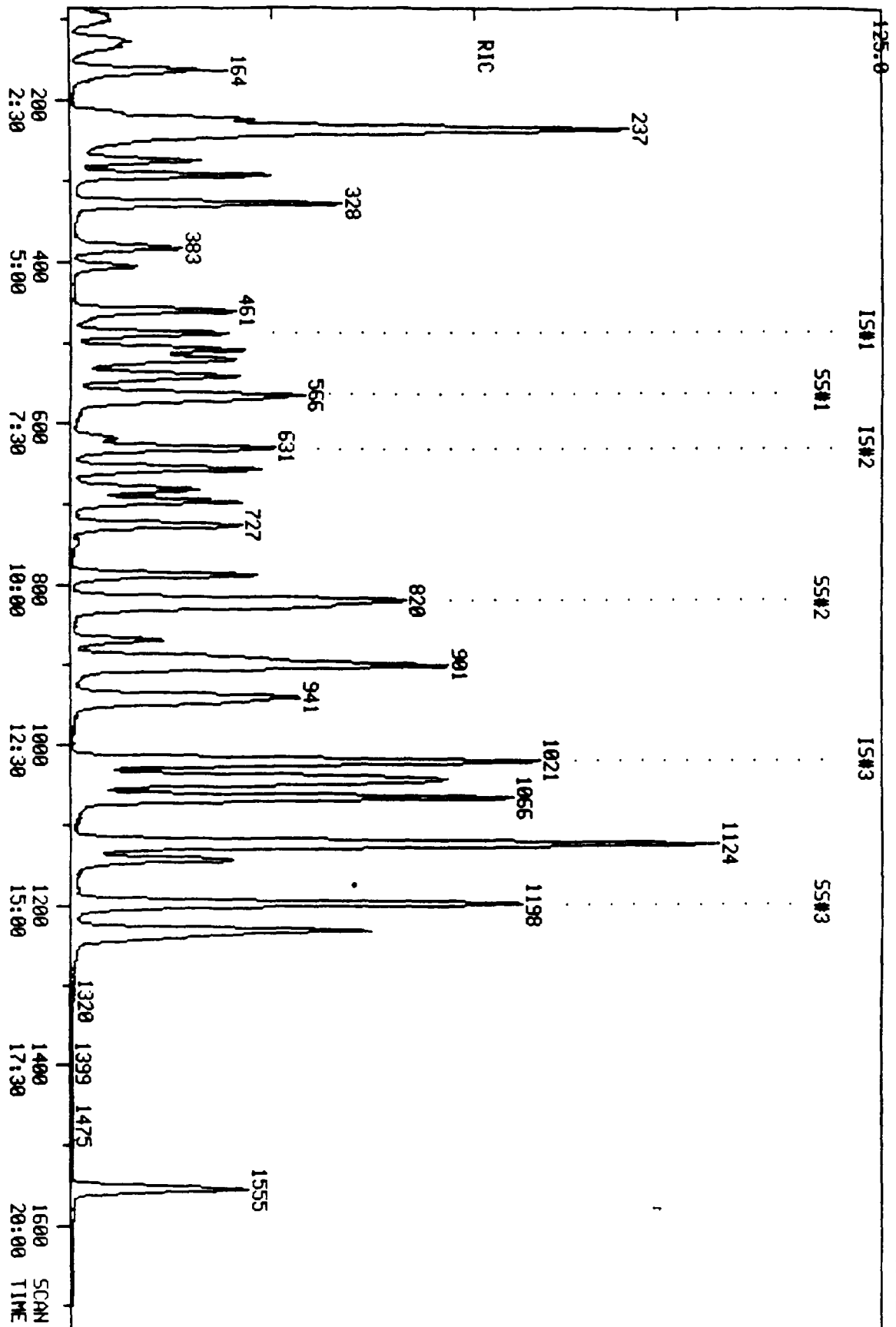
NO	REI(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	6:07	0.98	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:10	0.96	10.000	0.02	24.36	50.00	0.576	1.182	0.49
3	1:17	0.98	10.000	0.02	28.06	50.00	0.789	1.406	0.56
4	1:36	0.95	10.000	0.03	23.70	50.00	0.691	1.457	0.47
5	1:42	0.96	10.000	0.03	20.93	50.00	0.356	0.850	0.42
6	2:03	0.98	10.000	0.03	19.62	50.00	1.580	4.026	0.39
7	2:42	0.97	90.000	0.00	263.68	500.01	0.051	0.097	0.53
8	2:48	0.97	5.000	0.09	22.00	50.00	0.642	1.458	0.44
9	2:58	0.97	5.000	0.10	19.37	50.00	1.955	5.046	0.39
10	2:55	0.97	10.000	0.05	22.13	50.00	2.207	4.988	0.44
11	2:58	0.96	10.000	0.05	22.51	50.00	0.640	1.422	0.45
12	2:59	0.97	10.000	0.05	22.48	50.00	0.720	1.601	0.45
13	2:59	0.97	10.000	0.05	34.22	50.00	0.264	0.385	0.68
14	7:54	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
15	3:26	0.96	15.000	0.04	21.45	50.00	0.333	0.777	0.43
16	3:40	0.97	5.000	0.12	40.95	50.00	1.691	2.065	0.82
17	4:06	0.97	5.000	0.13	21.83	50.00	0.646	1.480	0.44
18	4:05	0.97	120.000	0.01	245.60	500.01	0.104	0.213	0.49
19	4:47	0.97	5.000	0.16	21.78	50.00	1.324	3.038	0.44
20	5:04	0.98	10.000	0.06	20.53	50.00	0.262	0.637	0.41
21	5:46	0.98	5.000	0.19	23.77	50.00	0.840	1.767	0.48
22	5:54	0.98	10.000	0.10	26.38	50.00	0.059	0.111	0.53
23	6:22	0.98	5.000	0.21	21.87	50.00	1.560	3.567	0.44
24	6:31	0.98	5.000	0.16	21.40	50.00	0.345	0.807	0.43
25	6:46	0.98	5.000	0.17	19.64	50.00	0.292	0.743	0.39
26	7:05	0.98	5.000	0.18	23.06	50.00	0.379	0.821	0.46
27	7:10	0.99	5.000	0.24	22.97	50.00	1.138	2.476	0.46
28	7:43	0.99	100.000	0.01	379.49	500.01	0.009	0.011	0.76
29	12:44	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
30	8:13	0.99	5.000	0.21	24.24	50.00	0.200	0.413	0.48
31	8:32	0.99	5.000	0.22	23.27	50.00	0.153	0.330	0.47
32	8:43	0.99	5.000	0.29	21.78	50.00	0.803	1.843	0.44
33	9:05	0.99	5.000	0.23	20.09	50.00	0.318	0.792	0.40
34	9:14		10.000			50.00		0.001	
35	9:52	0.99	5.000	0.25	22.31	50.00	0.354	0.793	0.45
36	10:15	0.99	15.000	0.05	24.32	50.00	0.163	0.336	0.49

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
37	10:22	0.99	5.000	0.16	22.43	50.00	0.295	0.657	0.45
38	10:52	0.99	5.000	0.28	22.24	50.00	0.172	0.386	0.44
39	11:07	0.99	5.000	0.28	24.37	50.00	0.169	0.346	0.49
40	11:13	0.99	10.000	0.09	22.78	50.00	0.251	0.550	0.46
41	11:16	0.99	5.000	0.18	23.03	50.00	0.237	0.514	0.46
42	11:43	0.99	15.000	0.06	26.23	50.00	0.118	0.225	0.52
43	11:45	1.00	5.000	0.30	21.51	50.00	0.278	0.647	0.43
44	11:50	1.00	5.000	0.30	21.44	50.00	0.288	0.671	0.43
45	12:47	1.00	5.000	0.20	23.09	50.00	0.425	0.921	0.46
46	13:00	1.00	5.000	0.33	20.61	50.00	0.273	0.662	0.41
47	13:05	1.00	5.000	0.21	22.10	50.00	0.193	0.436	0.44
48	13:19	1.00	5.000	0.21	20.43	50.00	0.350	0.857	0.41
49	14:01	1.00	5.000	0.22	20.40	50.00	0.292	0.716	0.41
50	14:04	1.00	5.000	0.22	21.61	50.00	0.516	1.194	0.43
51	14:17	1.00	5.000	0.37	21.64	50.00	0.244	0.564	0.43
52	14:57	1.00	15.000	0.13	20.11	50.00	0.087	0.217	0.40
53	15:22	1.00	15.000	0.08	21.67	50.00	0.081	0.188	0.43
54	15:25	1.00	5.000	0.24	19.88	50.00	0.220	0.553	0.40
55	15:31	1.00	15.000	0.08	19.54	50.00	0.064	0.164	0.39
56	19:26	1.00	10.000	0.15	48.05	100.00	0.085	0.176	0.48
57	7:03	0.98	5.000	0.23	20.76	50.00	1.080	2.601	0.42
58	14:58	1.00	5.000	0.24	21.41	50.00	0.424	0.991	0.43
59	10:15	0.99	5.000	0.16	22.71	50.00	0.508	1.119	0.45



RIC
 12/12/89 20:54:00
 SAMPLE: 10 ML EPA ID# USTD050 (STD# 1907) ON #19
 COND.S.:

COMPUTHER LABS
 COMPUTHER DATA: G1891212819 SCANS 86 TO 1700
 OUT OF 86 TO 1700
 217920.



QUANTITATION REPORT FILE: GT891212B19
DATA: GT891212B19.TI
12/12/89 20:54:00
SAMPLE: 10 ML EPA ID# VSTD050 (STD# 1907) DN #19
CONDS.:
SUBMITTED BY: 19 ANALYST: 1009

AMOUNT=AREA * REF. AMNT/(REF. AREA)* RESP. FACT)
RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1
2	221 CHLOROMETHANE <74-87-3> RO#2
3	231 VINYL CHLORIDE <75-01-4> RO#3
4	220 BROMOMETHANE <78-83-9> RO#4
5	209 CHLOROETHANE <75-00-3> RO#5
6	230 TRICHLOROFLUOROMETHANE <75-69-4> RO#6
7	201 ACROLEIN <107-02-8> RO#7
8	216 1,1-DICHLOROETHENE <75-35-4> RO#8
9	254 CARBON DISULFIDE <75-15-0> RO#9
10	285 IODOMETHANE <74-88-4> RO#10
11	297 1,1,1-TRICHLORO-2,2,2-TRIFLUOROETHANE <354-58-5> RO#11
12	266 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE <76-13-1> RO#12
13	252 ACETONE (2-PROPANONE) <67-64-1> RO#13
14	*249 1,4-DIFLUOROBENZENE (IS) <540-36-3> RO#14
15	298 3-CHLOROPROPENE <107-05-1> RO#15
16	222 METHYLENE CHLORIDE <75-09-2> RO#16
17	226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17
18	202 ACRYLONITRILE <107-13-1> RO#18
19	214 1,1-DICHLOROETHANE <75-34-3> RO#19
20	257 VINYL ACETATE <108-05-4> RO#20
21	237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21
22	253 2-BUTANONE <78-93-3> RO#22
23	211 CHLOROFORM <67-66-2> RO#23
24	227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24
25	206 CARBON TETRACHLORIDE <56-23-5> RO#25
26	203 BENZENE <71-43-2> RO#26
27	215 1,2-DICHLOROETHANE <107-06-2> RO#27
28	272 CROTONALDEHYDE <4170-30-3> RO#28
29	*270 D5-CHLOROENZENE (IS) RO#29
30	229 TRICHLOROETHENE <79-01-6> RO#30
31	217 1,2-DICHLOROPROPANE <78-87-5> RO#31
32	286 DIBROMOMETHANE <74-95-3> RO#32
33	212 BROMODICHLOROMETHANE <75-27-4> RO#33
34	210 2-CHLOROETHYL VINYL ETHER <110-75-8> RO#34
35	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35
36	256 4-METHYL-2-PENTANONE <108-01-1> RO#36
37	225 TOLUENE <108-88-3> RO#37
38	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38
39	228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39
40	287 ETHYLMETHACRYLATE <96-18-4> RO#40
41	224 TETRACHLOROETHENE <127-18-4> RO#41
42	255 2-HEXANONE <591-78-6> RO#42
43	208 DIBROMOCHLOROMETHANE <124-48-1> RO#43
44	245 1,2-DIBROMOMETHANE <1060-93-4> RO#44
45	207 CHLOROBENZENE <108-90-7> RO#45
46	273 1,1,1,2-TETRACHLOROETHANE <630-20-6> RO#46

NO NAME
 47 219 ETHYLBENZENE <100-41-4> RO#47
 48 330 M,P-XYLENE <133-02-7> RO#48
 49 239 O-XYLENE <133-02-7> RO#49
 50 251 STYRENE <100-42-5> RO#50
 51 205 BROMOFORM <75-25-2> RO#51
 52 274 CIS-1,4-DICHLORO-2-BUTENE <764-71-0> RO#52
 53 275 1,2,3-TRICHLOROPROPANE <96-18-4> RO#53
 54 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54
 55 290 TRANS-1,4-DICHLORO-2-BUTENE <110-57-6> RO#55
 56 262 1,2-DIBROMO-3-CHLOROPROPANE <96-12-8> RO#56
 57 #258 D4-1,2-DICHLOROETHANE RO#57
 58 #247 BROMOFLUOROBENZENE <460-00-4> RO#58
 59 #233 DB-TOLUENE RO#59

66-12-13-89
 12-13-89

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HQHT)	AMOUNT	%TOT
1	128	489	6:07	1	1.000	A BB	39810.	50.000 UG/KG	1.02
2	50	94	1:10	1	0.192	A BB	47052.	112.068 UG/KG	2.28
3	62	103	1:17	1	0.211	A BB	55979.	83.642 UG/KG	1.70
4	94	128	1:36	1	0.262	A BB	58007.	69.843 UG/KG	1.42
5	64	136	1:42	1	0.278	A BB	33828.	79.782 UG/KG	1.63
6	101	164	2:03	1	0.335	A BB	160277.	71.797 UG/KG	1.46
7	56	216	2:42	1	0.442	A BB	38751.	567.867 UG/KG	11.57
8	96	224	2:48	1	0.458	A BB	58056.	48.001 UG/KG	0.98
9	76	238	2:58	1	0.487	A BB	200896.	59.018 UG/KG	1.20
10	142	234	2:55	1	0.479	A BB	198558.	53.506 UG/KG	1.09
11	117	237	2:58	1	0.485	A BB	56602.	63.920 UG/KG	1.30
12	85	239	2:59	1	0.489	A BB	63752.	70.690 UG/KG	1.44
13	43	239	2:59	1	0.489	A BB	15339.	82.097 UG/KG	1.67
14	114	632	7:54	14	1.000	A BB	164799.	50.000 UG/KG	1.02
15	76	275	3:26	1	0.562	A BB	30919.	59.074 UG/KG	1.20
16	84	293	3:40	1	0.599	A BB	82213.	46.469 UG/KG	0.95
17	96	328	4:06	1	0.671	A BB	58926.	47.430 UG/KG	0.97
18	53	327	4:05	1	0.669	A BV	84666.	569.281 UG/KG	11.60
19	63	383	4:47	1	0.783	A BB	120952.	61.136 UG/KG	1.25
20	43	406	5:04	14	0.642	A BB	105000.	92.626 UG/KG	1.89
21	96	461	5:46	1	0.943	A BB	70339.	53.234 UG/KG	1.09
22	72	472	5:54	1	0.965	A VB	4435.	45.632 UG/KG	0.93
23	83	509	6:22	1	1.041	A BB	141995.	54.891 UG/KG	1.12
24	97	522	6:31	14	0.826	A BB	132958.	66.081 UG/KG	1.35
25	117	542	6:46	14	0.858	A BB	122382.	76.236 UG/KG	1.55
26	78	567	7:05	14	0.897	A BB	135273.	45.519 UG/KG	0.93
27	62	573	7:10	1	1.172	A BB	98568.	71.508 UG/KG	1.46
28	70	618	7:43	14	0.978	A BV	18606.	347.123 UG/KG	7.08
29	117	1019	12:44	29	1.000	A BB	191140.	50.000 UG/KG	1.02
30	130	657	8:13	14	1.040	A BB	68035.	42.286 UG/KG	0.86
31	63	683	8:32	14	1.081	A BB	54323.	55.361 UG/KG	1.13
32	174	697	8:43	1	1.425	A BB	73367.	51.312 UG/KG	1.05
33	83	727	9:05	14	1.150	A BB	130509.	60.688 UG/KG	1.24
34	63	NOT FOUND							
35	75	789	9:52	14	1.248	A BB	130611.	57.151 UG/KG	1.16
36	43	820	10:15	29	0.805	A BB	64211.	68.963 UG/KG	1.41
37	92	829	10:22	29	0.814	A BB	125522.	43.917 UG/KG	0.90
38	75	870	10:52	14	1.377	A BB	63543.	62.992 UG/KG	1.28
39	97	890	11:07	14	1.408	A BB	57072.	48.413 UG/KG	0.99
40	69	898	11:13	29	0.881	A BB	105155.	57.672 UG/KG	1.18
41	164	902	11:16	29	0.885	A BB	98327.	41.854 UG/KG	0.85

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HQHT)	AMOUNT	%TOT
42	43	938	11:43	29	0.921	A BB	42975.	66.919 UG/KG	1.36
43	129	940	11:45	14	1.487	A BB	106574.	64.078 UG/KG	1.31
44	107	947	11:50	14	1.498	A BB	110621.	62.679 UG/KG	1.28
45	112	1023	12:47	29	1.004	A BB	176126.	43.530 UG/KG	0.89
46	131	1040	13:00	14	1.646	A BB	109149.	72.672 UG/KG	1.48
47	106	1047	13:05	29	1.027	A BV	83326.	44.865 UG/KG	0.91
48	106	1066	13:19	29	1.046	A VB	163843.	53.586 UG/KG	1.09
49	106	1121	14:01	29	1.100	A BB	136822.	54.005 UG/KG	1.10
50	104	1125	14:04	29	1.104	A BB	228151.	52.441 UG/KG	1.07
51	173	1143	14:17	14	1.809	A BB	92870.	83.493 UG/KG	1.70
52	88	1196	14:57	14	1.892	A BB	35793.	84.106 UG/KG	1.71
53	110	1230	15:22	29	1.207	A BB	35883.	59.364 UG/KG	1.21
54	83	1233	15:25	29	1.210	A BB	105613.	42.187 UG/KG	0.86
55	53	1241	15:31	29	1.218	A BB	31321.	71.926 UG/KG	1.47
56	157	1555	19:26	29	1.526	A BB	67398.	110.003 UG/KG	2.24
57	65	564	7:03	1	1.153	A BB	103563.	85.063 UG/KG	1.73
58	95	1198	14:58	29	1.176	A BB	189399.	61.599 UG/KG	1.26
59	98	820	10:15	29	0.805	A BB	213794.	55.006 UG/KG	1.12

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	6:01	1.02	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:08	1.03	10.000	0.02	112.07	50.00	1.182	0.527	2.24
3	1:14	1.04	10.000	0.02	83.64	50.00	1.406	0.841	1.67
4	1:30	1.07	10.000	0.03	69.84	50.00	1.457	1.043	1.40
5	1:38	1.04	10.000	0.03	79.78	50.00	0.850	0.533	1.60
6	1:58	1.04	10.000	0.03	71.80	50.00	4.026	2.804	1.44
7	2:37	1.03	90.000	0.00	567.87	500.01	0.097	0.086	1.14
8	2:42	1.04	5.000	0.09	48.00	50.00	1.458	1.519	0.96
9	2:52	1.03	5.000	0.10	59.02	50.00	5.046	4.275	1.18
10	2:49	1.04	10.000	0.05	53.51	50.00	4.988	4.661	1.07
11	2:49	1.05	10.000	0.05	63.92	50.00	1.422	1.112	1.28
12	2:52	1.04	10.000	0.05	70.69	50.00	1.601	1.133	1.41
13	2:53	1.03	10.000	0.05	82.10	50.00	0.385	0.235	1.64
14	7:49	1.01	5.000	0.20	50.00	50.00	1.000	1.000	1.00
15	3:19	1.03	15.000	0.04	59.07	50.00	0.777	0.657	1.18
16	3:33	1.03	5.000	0.12	46.47	50.00	2.065	2.222	0.93
17	3:59	1.03	5.000	0.13	47.43	50.00	1.480	1.560	0.95
18	3:59	1.03	120.000	0.01	569.28	500.01	0.213	0.187	1.14
19	4:40	1.02	5.000	0.16	61.14	50.00	3.038	2.485	1.22
20	4:58	1.02	10.000	0.06	92.63	50.00	0.637	0.344	1.85
21	5:40	1.02	5.000	0.19	53.23	50.00	1.767	1.660	1.06
22	5:48	1.02	10.000	0.10	45.63	50.00	0.111	0.122	0.91
23	6:16	1.01	5.000	0.21	54.89	50.00	3.567	3.249	1.10
24	6:25	1.02	5.000	0.17	66.08	50.00	0.807	0.610	1.32
25	6:41	1.01	5.000	0.17	76.24	50.00	0.743	0.487	1.52
26	7:01	1.01	5.000	0.18	45.52	50.00	0.821	0.902	0.91
27	7:05	1.01	5.000	0.23	71.51	50.00	2.476	1.731	1.43
28	7:40	1.01	100.000	0.01	347.12	500.01	0.011	0.016	0.69
29	12:42	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
30	8:09	1.01	5.000	0.21	42.29	50.00	0.413	0.488	0.85
31	8:28	1.01	5.000	0.22	55.36	50.00	0.330	0.298	1.11
32	8:39	1.01	5.000	0.29	51.31	50.00	1.843	1.796	1.03
33	9:02	1.01	5.000	0.23	60.69	50.00	0.792	0.652	1.21
34	9:14		10.000			50.06		0.001	
35	9:49	1.01	5.000	0.25	57.15	50.00	0.793	0.693	1.14
36	10:12	1.00	15.000	0.05	68.96	50.00	0.336	0.244	1.38

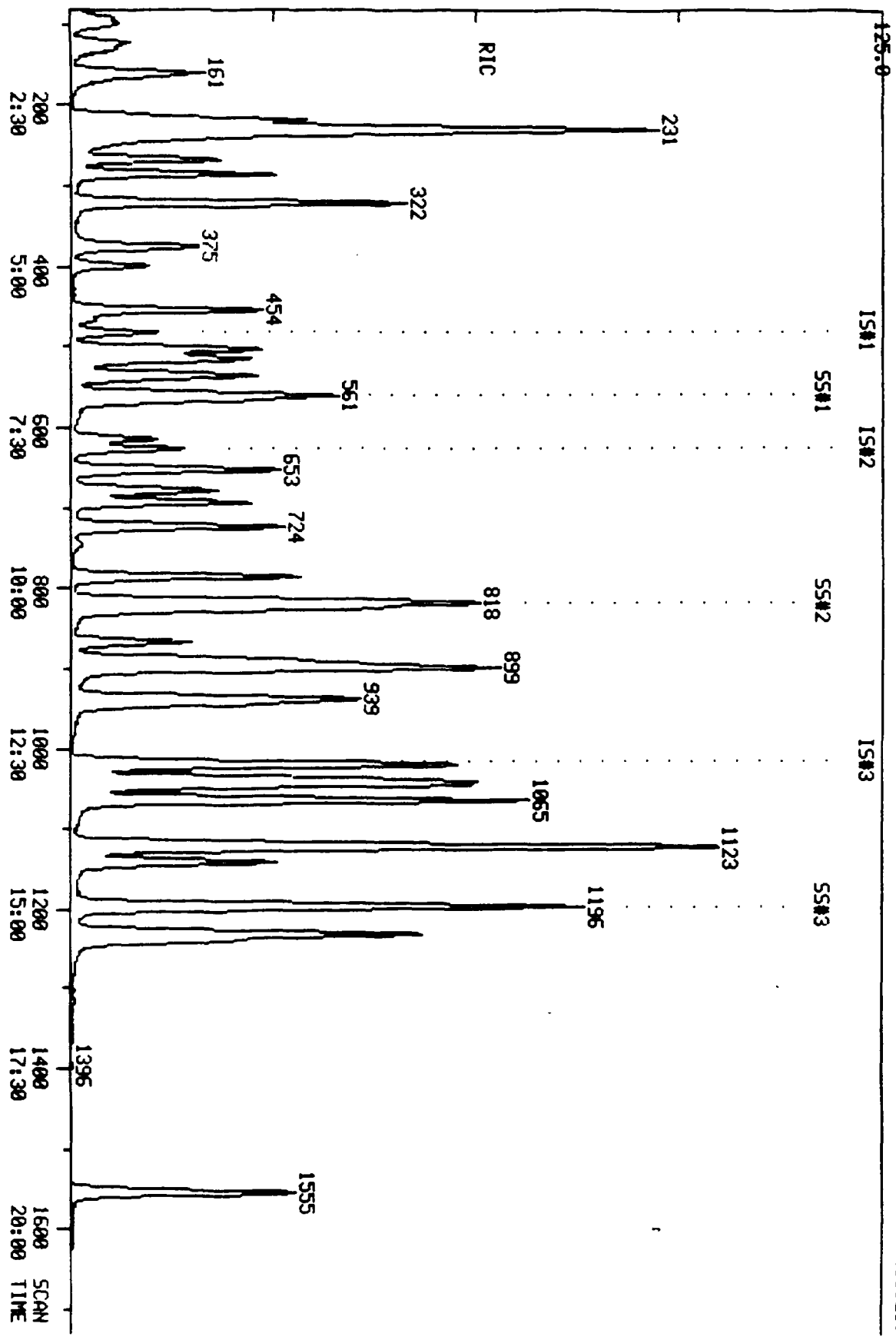
NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
37	10:18	1.01	5.000	0.16	43.92	50.00	0.657	0.748	0.88
38	10:49	1.00	5.000	0.28	62.99	50.00	0.386	0.306	1.26
39	11:05	1.00	5.000	0.28	48.41	50.00	0.346	0.358	0.97
40	11:10	1.00	10.000	0.09	57.67	50.00	0.550	0.477	1.15
41	11:14	1.00	5.000	0.18	41.85	50.00	0.514	0.615	0.84
42	11:41	1.00	15.000	0.06	66.92	50.00	0.225	0.168	1.34
43	11:42	1.00	5.000	0.30	64.08	50.00	0.647	0.505	1.28
44	11:48	1.00	5.000	0.30	62.68	50.00	0.671	0.535	1.25
45	12:46	1.00	5.000	0.20	43.53	50.00	0.921	1.058	0.87
46	12:58	1.00	5.000	0.33	72.67	50.00	0.662	0.456	1.45
47	13:03	1.00	5.000	0.21	44.87	50.00	0.436	0.486	0.90
48	13:17	1.00	5.000	0.21	53.59	50.00	0.857	0.800	1.07
49	13:59	1.00	5.000	0.22	54.00	50.00	0.716	0.663	1.08
50	14:02	1.00	5.000	0.22	52.44	50.00	1.194	1.138	1.05
51	14:15	1.00	5.000	0.36	83.49	50.00	0.564	0.337	1.67
52	14:55	1.00	15.000	0.13	84.11	50.00	0.217	0.129	1.68
53	15:22	1.00	15.000	0.08	59.36	50.00	0.188	0.158	1.19
54	15:23	1.00	5.000	0.24	42.19	50.00	0.553	0.655	0.84
55	15:30	1.00	15.000	0.08	71.93	50.00	0.164	0.114	1.44
56	19:25	1.00	10.000	0.15	110.00	100.00	0.176	0.160	1.10
57	6:58	1.01	5.000	0.23	85.06	50.00	2.601	1.529	1.70
58	14:57	1.00	5.000	0.24	61.60	50.00	0.991	0.804	1.23
59	10:13	1.00	5.000	0.16	55.01	50.00	1.119	1.017	1.10

RIC
 12/12/89 23:47:00
 SAMPLE: 10ML EPA ID# USTD100 (STD# 1908) ON #19
 COND.:

COMPUCHEM LABS

COMPUCHEM DATA: GM891212B19 SCANS 82 TO 1700
 OUT OF 82 TO 1700

405120.



QUANTITATION REPORT FILE: GW891212B19
DATA: GW891212B19.TI
12/12/89 23:47:00
SAMPLE: 10ML EPA ID# VSTD100 (STD# 1908) ON #19
CONDS.:
SUBMITTED BY: 19 ANALYST: 1009

AMOUNT=AREA * REF.AMNT/(REF.AREA)* RESP.FACT)
RESP. FAC. FROM LIBRARY ENTRY

NO	NAME
1	*234 BROMOCHLOROMETHANE (IS) <75-97-5> RO#1
2	221 CHLOROMETHANE <74-87-3> RO#2
3	231 VINYL CHLORIDE <75-01-4> RO#3
4	220 BROMOMETHANE <78-83-9> RO#4
5	209 CHLOROETHANE <75-00-3> RO#5
6	230 TRICHLOROFLUOROMETHANE <75-69-4> RO#6
7	201 ACRYLEIN <107-02-8> RO#7
8	216 1,1-DICHLOROETHENE <75-35-4> RO#8
9	254 CARBON DISULFIDE <75-15-0> RO#9
10	285 IODOMETHANE <74-88-4> RO#10
11	297 1,1,1-TRICHLORO-2,2,2-TRIFLUOROETHANE <354-58-5> RO#11
12	266 1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE <76-13-1> RO#12
13	252 ACETONE (2-PROPANONE) <67-64-1> RO#13
14	*248 1,4-DIFLUOROBENZENE (IS) <540-36-3> RO#14
15	298 3-CHLOROPROPENE <107-05-1> RO#15
16	222 METHYLENE CHLORIDE <75-09-2> RO#16
17	226 TRANS-1,2-DICHLOROETHENE <156-60-5> RO#17
18	202 ACRYLONITRILE <107-13-1> RO#18
19	214 1,1-DICHLOROETHANE <75-34-3> RO#19
20	257 VINYL ACETATE <108-05-4> RO#20
21	237 CIS-1,2-DICHLOROETHENE <156-59-2> RO#21
22	253 2-BUTANONE <78-93-3> RO#22
23	211 CHLOROFORM <67-66-2> RO#23
24	227 1,1,1-TRICHLOROETHANE <71-55-6> RO#24
25	206 CARBON TETRACHLORIDE <56-23-5> RO#25
26	203 BENZENE <71-43-2> RO#26
27	215 1,2-DICHLOROETHANE <107-06-2> RO#27
28	272 CROTONALDEHYDE <4170-30-3> RO#28
29	*270 D5-CHLOROBENZENE (IS) RO#29
30	229 TRICHLOROETHENE <79-01-6> RO#30
31	217 1,2-DICHLOROPROPANE <78-87-5> RO#31
32	286 DIBROMOMETHANE <74-95-3> RO#32
33	212 BROMODICHLOROMETHANE <75-27-4> RO#33
34	210 2-CHLOROETHYL VINYL ETHER <110-75-8> RO#34
35	218 CIS-1,3-DICHLOROPROPENE <10061-1-5> RO#35
36	256 4-METHYL-2-PENTANONE <108-01-1> RO#36
37	225 TOLUENE <108-88-3> RO#37
38	250 TRANS-1,3-DICHLOROPROPENE <10061-02-6> RO#38
39	228 1,1,2-TRICHLOROETHANE <79-00-5> RO#39
40	287 ETHYLMETHACRYLATE <96-18-4> RO#40
41	224 TETRACHLOROETHENE <127-18-4> RO#41
42	255 2-HEXANONE <591-78-6> RO#42
43	208 DIBROMOCHLOROMETHANE <124-48-1> RO#43
44	245 1,2-DIBROMOMETHANE <1060-93-4> RO#44
45	207 CHLOROBENZENE <108-90-7> RO#45
46	273 1,1,1,2-TETRACHLOROETHANE <630-20-6> RO#46

NO NAME
 47 219 ETHYLBENZENE <100-41-4> RO#47
 48 330 M,P-XYLENE <133-02-7> RO#48
 49 239 O-XYLENE <133-02-7> RO#49
 50 251 STYRENE <100-42-5> RO#50
 51 205 BROMOFORM <75-25-2> RO#51
 52 274 CIS-1,4-DICHLORO-2-BUTENE <764-71-0> RO#52
 53 275 1,2,3-TRICHLOROPROPANE <96-18-4> RO#53
 54 223 1,1,2,2-TETRACHLOROETHANE <79-34-5> RO#54
 55 290 TRANS-1,4-DICHLORO-2-BUTENE <110-57-6> RO#55
 56 262 1,2-DIBROMO-3-CHLOROPROPANE <96-12-8> RO#56
 57 #258 D4-1,2-DICHLOROETHANE RO#57
 58 #247 BROMOFLUOROBENZENE <460-00-4> RO#58
 59 #233 DB-TOLUENE RO#59

OKS
 12.13.89
 16.11.89

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	483	6:02	1	1.000	A BB	40867.	50.000 UG/KG	0.46
2	50	92	1:09	1	0.190	A BB	102274.	105.871 UG/KG	0.98
3	62	100	1:15	1	0.207	A BB	124707.	108.506 UG/KG	1.01
4	94	122	1:31	1	0.253	A BB	105280.	88.401 UG/KG	0.82
5	64	135	1:41	1	0.280	A BB	52430.	75.490 UG/KG	0.70
6	101	161	2:01	1	0.333	A BB	280196.	85.149 UG/KG	0.79
7	56	214	2:40	1	0.443	A BB	105028.	1320.120 UG/KG	12.27
8	96	219	2:44	1	0.453	A BB	123305.	103.448 UG/KG	0.96
9	76	232	2:54	1	0.480	A BB	361519.	87.649 UG/KG	0.81
10	142	229	2:52	1	0.474	A BB	371437.	91.114 UG/KG	0.85
11	117	230	2:52	1	0.476	A BB	109646.	94.352 UG/KG	0.88
12	85	234	2:55	1	0.484	A BB	128084.	97.857 UG/KG	0.91
13	43	236	2:57	1	0.489	A BB	29166.	92.609 UG/KG	0.86
14	114	627	7:50	14	1.000	A BB	160552.	50.000 UG/KG	0.46
15	76	268	3:21	1	0.555	A BB	69198.	109.008 UG/KG	1.01
16	84	286	3:34	1	0.592	A BB	157293.	93.187 UG/KG	0.87
17	96	320	4:00	1	0.663	A BB	127522.	105.407 UG/KG	0.98
18	53	322	4:01	1	0.667	A BB	231597.	1332.360 UG/KG	12.39
19	63	375	4:41	1	0.776	A BB	275599.	110.982 UG/KG	1.03
20	43	399	4:59	14	0.636	A BB	239151.	116.895 UG/KG	1.09
21	96	455	5:41	1	0.942	A BB	153530.	106.313 UG/KG	0.99
22	72	467	5:50	1	0.967	A VB	12310.	135.193 UG/KG	1.26
23	83	504	6:18	1	1.043	A BB	308267.	105.741 UG/KG	0.98
24	97	515	6:26	14	0.821	A BB	283449.	109.413 UG/KG	1.02
25	117	536	6:42	14	0.855	A BB	261110.	109.500 UG/KG	1.02
26	78	562	7:01	14	0.896	A BB	291212.	110.486 UG/KG	1.03
27	62	569	7:07	1	1.178	A BB	215786.	106.629 UG/KG	0.99
28	70	614	7:40	14	0.979	A BV	75840.	2092.000 UG/KG	19.45
29	117	1017	12:43	29	1.000	A BB	179393.	50.000 UG/KG	0.46
30	130	653	8:10	14	1.041	A BB	149206.	112.555 UG/KG	1.05
31	63	679	8:29	14	1.083	A BB	115808.	109.412 UG/KG	1.02
32	174	694	8:40	1	1.437	A BB	152374.	101.158 UG/KG	0.94
33	83	724	9:03	14	1.155	A BB	298541.	117.401 UG/KG	1.09
34	63	NOT FOUND							
35	75	786	9:49	14	1.254	A BB	310849.	122.146 UG/KG	1.14
36	43	818	10:13	29	0.804	A BB	161789.	134.231 UG/KG	1.25
37	92	827	10:20	29	0.813	A BB	272759.	115.764 UG/KG	1.08
38	75	868	10:51	14	1.384	A BB	152826.	123.435 UG/KG	1.15
39	97	888	11:06	14	1.416	A BB	135450.	121.805 UG/KG	1.13
40	69	895	11:11	29	0.880	A BB	239580.	121.377 UG/KG	1.13
41	164	900	11:15	29	0.885	A BB	202584.	109.761 UG/KG	1.02

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
42	43	936	11:42	29	0.920	A BB	138119.	171.217 UG/KG	1.59
43	129	939	11:44	14	1.498	A BB	236586.	113.933 UG/KG	1.06
44	107	945	11:49	14	1.507	A BB	231901.	107.591 UG/KG	1.00
45	112	1021	12:46	29	1.004	A BB	377723.	114.252 UG/KG	1.06
46	131	1039	12:59	14	1.657	A BB	215170.	101.175 UG/KG	0.94
47	106	1046	13:04	29	1.029	A BV	176083.	112.577 UG/KG	1.05
48	105	1065	13:19	29	1.047	A VB	307441.	99.965 UG/KG	0.93
49	106	1120	14:00	29	1.101	A BB	253339.	98.642 UG/KG	0.92
50	104	1124	14:03	29	1.105	A BB	434682.	101.500 UG/KG	0.94
51	173	1141	14:16	14	1.820	A BB	219162.	121.116 UG/KG	1.13
52	88	1195	14:56	14	1.906	A BB	85401.	122.454 UG/KG	1.14
53	110	1229	15:22	29	1.208	A BB	77032.	114.366 UG/KG	1.06
54	83	1232	15:24	29	1.211	A BB	248074.	125.135 UG/KG	1.16
55	53	1241	15:31	29	1.220	A BB	74508.	126.729 UG/KG	1.18
56	157	1554	19:25	29	1.528	A BB	165105.	261.012 UG/KG	2.43
57	65	559	6:59	1	1.157	A BB	223256.	105.000 UG/KG	0.98
58	95	1197	14:58	29	1.177	A BB	395588.	111.271 UG/KG	1.03
59	98	817	10:13	29	0.803	A BB	463188.	115.419 UG/KG	1.07

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
1	6:07	0.99	10.000	0.10	50.00	50.00	1.000	1.000	1.00
2	1:10	0.98	10.000	0.02	105.87	50.00	2.503	1.182	2.12
3	1:17	0.97	10.000	0.02	108.51	50.00	3.052	1.406	2.17
4	1:36	0.95	10.000	0.03	88.40	50.00	2.576	1.457	1.77
5	1:42	0.99	10.000	0.03	75.49	50.00	1.283	0.850	1.51
6	2:03	0.98	10.000	0.03	85.15	50.00	6.856	4.026	1.70
7	2:42	0.99	90.000	0.00	1320.13	500.01	0.257	0.097	2.64
8	2:48	0.98	5.000	0.09	103.45	50.00	3.017	1.458	2.07
9	2:58	0.97	5.000	0.10	87.65	50.00	8.846	5.046	1.75
10	2:55	0.98	10.000	0.05	91.11	50.00	9.089	4.988	1.82
11	2:58	0.97	10.000	0.05	94.35	50.00	2.683	1.422	1.89
12	2:59	0.98	10.000	0.05	97.86	50.00	3.134	1.601	1.96
13	2:59	0.99	10.000	0.05	92.61	50.00	0.714	0.385	1.85
14	7:54	0.99	5.000	0.20	50.00	50.00	1.000	1.000	1.00
15	3:26	0.97	15.000	0.04	109.01	50.00	1.693	0.777	2.18
16	3:40	0.98	5.000	0.12	93.19	50.00	3.849	2.065	1.86
17	4:06	0.98	5.000	0.13	105.41	50.00	3.120	1.480	2.11
18	4:05	0.98	120.000	0.01	1332.36	500.01	0.567	0.213	2.66
19	4:47	0.98	5.000	0.16	110.98	50.00	6.744	3.038	2.22
20	5:04	0.98	10.000	0.06	116.89	50.00	1.490	0.637	2.34
21	5:46	0.99	5.000	0.19	106.31	50.00	3.757	1.767	2.13
22	5:54	0.99	10.000	0.10	135.19	50.00	0.301	0.111	2.70
23	6:22	0.99	5.000	0.21	105.74	50.00	7.543	3.567	2.11
24	6:31	0.99	5.000	0.16	109.41	50.00	1.765	0.807	2.19
25	6:46	0.99	5.000	0.17	109.50	50.00	1.626	0.743	2.19
26	7:05	0.99	5.000	0.18	110.49	50.00	1.814	0.821	2.21
27	7:10	0.99	5.000	0.24	106.63	50.00	5.280	2.476	2.13
28	7:43	0.99	100.000	0.01	2092.00	500.01	0.047	0.011	4.18
29	12:44	1.00	5.000	0.20	50.00	50.00	1.000	1.000	1.00
30	8:13	0.99	5.000	0.21	112.55	50.00	0.929	0.413	2.25
31	8:32	0.99	5.000	0.22	109.41	50.00	0.721	0.330	2.19
32	8:43	1.00	5.000	0.29	101.16	50.00	3.729	1.843	2.02
33	9:05	1.00	5.000	0.23	117.40	50.00	1.859	0.792	2.35
34	9:14		10.000			50.00		0.001	
35	9:52	1.00	5.000	0.25	122.15	50.00	1.936	0.793	2.44
36	10:15	1.00	15.000	0.05	134.23	50.00	0.902	0.336	2.68

NO	RET(L)	RATIO	RRT(L)	RATIO	AMNT	AMNT(L)	R. FAC	R. FAC(L)	RATIO
37	10:22	1.00	5.000	0.16	115.76	50.00	1.520	0.657	2.32
38	10:52	1.00	5.000	0.28	123.43	50.00	0.952	0.386	2.47
39	11:07	1.00	5.000	0.28	121.81	50.00	0.844	0.346	2.44
40	11:13	1.00	10.000	0.09	121.38	50.00	1.336	0.950	2.43
41	11:16	1.00	5.000	0.18	109.76	50.00	1.129	0.514	2.20
42	11:43	1.00	15.000	0.06	171.22	50.00	0.770	0.225	3.42
43	11:45	1.00	5.000	0.30	113.93	50.00	1.474	0.647	2.28
44	11:50	1.00	5.000	0.30	107.59	50.00	1.444	0.671	2.15
45	12:47	1.00	5.000	0.20	114.25	50.00	2.106	0.921	2.29
46	13:00	1.00	5.000	0.33	101.17	50.00	1.340	0.662	2.02
47	13:05	1.00	5.000	0.21	112.58	50.00	0.982	0.436	2.25
48	13:19	1.00	5.000	0.21	99.97	50.00	1.714	0.857	2.00
49	14:01	1.00	5.000	0.22	98.64	50.00	1.412	0.716	1.97
50	14:04	1.00	5.000	0.22	101.50	50.00	2.423	1.194	2.03
51	14:17	1.00	5.000	0.36	121.12	50.00	1.365	0.564	2.42
52	14:57	1.00	15.000	0.13	122.45	50.00	0.532	0.217	2.45
53	15:22	1.00	15.000	0.08	114.37	50.00	0.429	0.188	2.29
54	15:25	1.00	5.000	0.24	125.14	50.00	1.383	0.553	2.50
55	15:31	1.00	15.000	0.08	126.73	50.00	0.415	0.164	2.53
56	19:26	1.00	10.000	0.15	261.01	100.00	0.460	0.176	2.61
57	7:03	0.99	5.000	0.23	105.00	50.00	5.463	2.601	2.10
58	14:58	1.00	5.000	0.24	111.27	50.00	2.205	0.991	2.23
59	10:15	1.00	5.000	0.16	115.42	50.00	2.582	1.119	2.31