



5 Underwood Ct., Delran, NJ 08075 609-461-4003

1S8112

TECHNICAL ASSISTANCE TEAM FOR EMERGENCY RESPONSE REMOVAL AND PREVENTION
EPA CONTRACT 68-WO-0036

MEMORANDUM

TO: Jamie Fenske, OSC, EPA, Region III
Western Response Section TDD #9111-26
PCS #2035

THRU: Mike Zickler, TATL, Region III M⁷

FROM: Marian Murphy, TAT Region III M¹¹

SUBJECT: Big John's Samples Analytical Review

DATE: January 10, 1992

This report covers the review of the analytical data package for five (5) drum samples and one (1) solid sample collected at the Big John's Site on December 10, 1991. The samples were received at Princeton Testing on December 12, 1991. Three drums were analyzed for volatile organics (VOA) and flashpoint. One drum was analyzed for PCBs, and one drum was analyzed for sulfites. The solid sample was analyzed for asbestos.

ANALYTICAL METHODOLOGY

The VOA samples were analyzed in accordance with EPA Method 8240. The PCB sample was analyzed in accordance with EPA Method 8080. The flashpoint samples were analyzed in accordance with EPA Method 1020. The sulfite analysis was performed in accordance with EPA Method 377.1. The solid was analyzed for asbestos by PLM.

- Signed chain-of-custody records were received.
- The GC/MS tune data for the VOAs and the internal standard data were acceptable. The initial and continuing calibration data had some compounds out of QC criteria, however, only 2-butanone was qualified approximate, for sample D01. The matrix spike/matrix spike duplicate recoveries, surrogate spike recoveries and relative percent difference (RPD) values were acceptable for the most part. The method blank contained methylene chloride and acetone. Samples D01, D05 and D06 should be considered not detected for methylene chloride, and sample D06 should be considered not detected for acetone.

Roy E. Weston, Inc.

MAJOR PROGRAMS DIVISION

In Association with Foster Wheeler Enviresponse, Inc., Resource Applications, C.C. Johnson & Malhotra, P.C., R.E. Sarriera Associates, and GRB Environmental Services, Inc.

MR100629

Big John's Samples Analytical Review
January 10, 1992
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- The PCB calibration data was acceptable. The method blank was free of contamination. The surrogate spike recovery for sample D04 was lower than the QC limits, but no PCBs were found, therefore no data was qualified. As no PCBs were found, no second column confirmations were performed.

CONCLUSION

Accept the data as presented except for methylene chloride, acetone and 2-butanone. The methylene chloride should be considered not detected for samples D01, D05 and D06 due to blank contamination. The acetone should be considered not detected for sample D06 due to blank contamination. The 2-butanone should be considered approximate for sample D01 due to calibration outliers. It should be noted that any volatile analysis marked "E" should be considered approximate since it is an estimated concentration.

MM/mr

AR100630

Attn: Mr. John J. Weitzel
Curtis Ridge, Suite 1000
Philadelphia, Pennsylvania 19106

CHAIN OF CUSTODY RECORD

PROJ. NO.	PROJECT NAME	NO. OF CONTAINERS	REMARKS
	Ba. Toh's Garage		
SAMPLERS (Signature)			
<i>John Michael P. Jones</i>			
STA. NO.	DATE	TIME	STATION LOCATION
D-21	12/10	04:22	X Inside Garage
D-22	12/10	04:22	X Inside Garage
D-23	12/10	04:22	X Inside Garage
D-24	12/10	04:22	X Inside Garage
D-25	12/10	04:22	X Inside Garage
D-26	12/10	04:22	X Inside Garage
D-27	12/10	04:22	X Inside Garage
			Please send one copy of Report to: PAT Battaglia 1/2 Pay F. Weitzel, 141 W. 2nd St., Philadelphia, PA 19103 One copy to Maria A. Murphy, the Inspector.
			5 Landenbach Court Delran, NJ 07075
Reinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time Received by: (Signature)
<i>John Michael P. Jones</i>	12/10/82 04:22	<i>John Michael P. Jones</i>	
Reinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time Received by: (Signature)
<i>John Michael P. Jones</i>			
Reinquished by: (Signature)	Date / Time	Received for Laboratory by: (Signature)	Date / Time Remarks
			<i>✓</i>

• Principals to DMR
Secretary for

Weston TAT
5 Underwood Court
Delran, New Jersey 08075
Attn: Marian Murphy

JOB # 9108360-001 / 2GMWH

AR100632



Weston TAT
5 Underwood Court
Delran, New Jersey 08075
Attn: Marian Murphy

January 3, 1992
Job #: 9108360-001/2GMWH

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AR100633*

APPROVED SAMPLE ANALYSIS REQUEST

Weston TAT
5 Underwood Court
Delran, New Jersey 08075

Attention: Marian Murphy
Phone: (609) 461-4003

Customer Number: T1342
Credit By: Heather Menzel
Approved By: Susan Sharples

Project No.: 9108360-001GMNH

Date Received: 12/13/91
Analysis Due : 12/25/91

Number Of Samples : 6
Number Of Containers: 6

Reports: Standard Report Format

Sample I.D.'s P.T.L.--Client	Requested Analytical Services
001 -- Sample D01 -----	Flashpoint Purgeable Volatile Organics By Method 8240 Page 1 Purgeable Volatile Organics By Method 8240 Page 2
002 -- Sample D03 -----	Sulfite (As SO3) In Solid Phase
003 -- Sample D04 -----	Polychlorinated Biphenyls In Oil
004 -- Sample D05 -----	Flashpoint Volatile Organics By GCMS -- Method 624 Page 1 Volatile Organics By GCMS -- Method 624 Page 2
005 -- Sample D06 -----	Flashpoint Volatile Organics By GCMS -- Method 624 Page 1 Volatile Organics By GCMS -- Method 624 Page 2
006 -- Sample A07 -----	Asbestos In Bulk/Non-Aqueous Sample By PLM

Notes:

1. Samples originated from Weston, Wheeling, WV 26003.
2. Extra copy of report to: Roy F. Weston, Inc.
141 Waddles Run Rd., Wheeling WV 26003.
Attn: Pat Bishop
3. Samples came in after 11:00 a.m. 12/11/91.
4. Per Client's instructions to H Menzel for Sample D04, run only the oil portion for PCB. Flashpoint to All Service.
5. Please see attached list for Detection limit.
6. NOTE THAT FOR RESULTS EXCEEDING THE PRESCRIBED TURNAROUND TIME (IN CALENDAR DAYS), A REDUCTION IN THE TOTAL CHARGE WILL BE THREE PERCENT FOR THE FIRST DAY, TWO PERCENT FOR THE SECOND DAY, AND AN ADDITIONAL ONE PERCENT FOR EACH ADDITIONAL DAY.

Job Type: Standard/PLM

Received By Lab:	Initials/Date
Reviewed By:	
Q.A. Approved:	

Printed By: Chris Jamieson
Date: 01/07/92
AR | O'Brien 16:57:12

DRAFT

DRAFT

CHAIN OF CUSTODY RECORD

CLIENT <i>Mobil</i>	SAMPLERS							PTL. JOB NO. <i>741-304</i>
CONTACT <i>A. G. H.</i>	FIELD SUPERVISOR							CLIENT PROJECT NO.
PHONE NO.	COMPANY							Page _____ of _____
SAMPLING SITE	PHONE NO.							REMARKS (Specify QA/QC, preservation required, due date, etc.)
SAMPLE IDENTIFICATION	MATRIX	COLLECTION		CONTAINER			ANALYSES REQUIRED (Specify Method if Known)	
		COMP.	GAS	DATE	TIME	TYPE	NO.	
<i>No 1</i>	<i>Gas</i>			<i>G</i>			<i>Flushed, 6246</i>	
<i>No 2</i>	<i>Gas</i>			<i>G</i>			<i>Supply</i>	
<i>No 5</i>	<i>Gas</i>			<i>G</i>			<i>Flashed, 624</i>	
<i>No 6</i>	<i>Gas</i>			<i>G</i>			<i>Flashed, 624</i>	
<i>No 7</i>	<i>Gas</i>			<i>G</i>			<i>Supply by PTL</i>	
LABORATORY SEAL NO. <i>C5</i>	LABORATORY SEAL NO. _____							FIELD SEAL BROKEN BY _____
SEALED ON _____ <i>12/16/99</i>	AT _____ <i>Military Time:</i> <i>10:00 AM</i>	UNSEALED ON _____ <i>Date:</i> <i>12/16/99</i>	AT _____ <i>Military Time:</i> <i>10:00 AM</i>	SAMPLING COMPLETED ON _____ <i>Date:</i> <i>12/16/99</i>	AT _____ <i>Military Time:</i> <i>10:00 AM</i>	FIELD SEAL NO. <i>1730</i>	DATE _____ <i>12/16/99</i>	TIME _____ <i>10:00 AM</i>
PREPARED BY _____			RECEIVED BY (Print and Sign Name): <i>Mark V. T. / V. T.</i>			REMARKS _____	REASON FOR CHANGE OF CUSTODY <i>1. Not at site</i>	
MEANS OF TRANSPORT _____								
RELINGUISHED BY (Print and Sign Name): <i>J. B. D. / J. B. D.</i>								

princeton testing
laboratory inc.

Weston TAT
5 Underwood Court
Delran, New Jersey 08075
Attention: Marian Murphy

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-6
FAX (609) 452-6

Report Date: 01/02/92
Job Number: 9108360-001GMWH
Date Received: 12/13/91

Analyses	Sample I.D.:	Sample D01	Sample D03	Sample D05
Flashpoint in degrees fahrenheit		<50° F Flash	N/R 67,800	<50° F N/R
Sulfite in mg/kg		N/R	67,800	67,7%

Analysis	Sample I.D.:	Sample D06
Flashpoint in degrees fahrenheit		139° F + Combustible

N/R = Not Requested

Eduardo A. Almeida
Eduardo A. Almeida, V.P., Manager
Water, Wastewater, and Microbiology

AB 100636

For inquiries call us at (609) 452-9050 and ask for our Customer Service Department

Member: American Council of Independent Laboratories, Inc.

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princeton testing laboratory inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
FAX (609) 452-0347

Weston TAT
5 Underwood Court
Delran, New Jersey 08075
Attention: Marian Murphy

Report Date: 01/07/92
Job Number: 9108360-001CMWH
Date Received: 12/13/91

Analysis: Polychlorinated Biphenyls In Oil
Units: mg/kg

Parameters	Sample I.D.	Sample D04	Blank 12/23/91
Aroclor 1221		<5	<5
Aroclor 1232		<5	<5
Aroclor 1242		<5	<5
Aroclor 1248		<5	<5
Aroclor 1254		<5	<5
Aroclor 1260		<5	<5
Aroclor 1016		<5	<5

RECOVERY DATA

* Aldrin (Surr.) (55-1458) 478 1598

NOTE: Low recovery due to extensive sulfuric acid cleanup.

Jane Dennison, Supervisor
Gas Chromatography Laboratory

For inquiries call us at (609) 452-9050  Customer Service Department

Member: American Council of Independent Laboratories, Inc.

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Software Version: 3.01.02

Date: 1/23/92 10:57

Sample Name: 1560-1000-100

Data File: C:\V2000\04740.1217027.D, Date: 10-16-91, File:

Sequence File: none, Cycle: 25, Standard: A

Instrument: VARIAN 3400, Recorder: 1, Detector: 10000,

Sample Amount: 1.0000, Dilution Factor: 1.00

Princeton Testing Laboratory Analysis Report

PESTICIDES AND PCB'S BY CAPILLARY CHROMATOGRAPHY

Peak Ret Time # [min]	Component Name	Key Agent Concentration [ppm]	[ppb]	Area	Weight in Microg [g]	Reference	DL, Detection Limit	Dil. Dilution Factor
1	0.140	-0.1000	-0.1000	18560.00	1.8560E-06	-----	-0.1000	
2	0.150	1.0000	1.0000	235040.00	2.3504E-05	-----	1.0000	
3	0.410	0.7000	0.7000	766000.00	7.6600E-05	-----	1.0000	
4	6.500 D-800	0.1000	0.1000	196701.00	1.9670E-06	-0.0010 DL	-0.0001	
5	8.610 HEPTACHLOR	-0.1000	-0.1000	110501.00	1.1050E-06	0.1105 DL	-0.1000	
6	8.684	-0.1000	-0.1000	130401.00	1.3040E-06	-----	-0.1000	
7	9.500 D-400	0.1000	0.1000	407450.00	4.0745E-06	0.1000 DL	0.1000	
8	10.044 HEPTACHLOR EPIC	0.1000	0.1000	168046.00	1.6804E-06	-0.1000 DL	0.1000	
9	12.045 DIBENZIN	0.1000	0.1000	801000.00	8.0100E-06	-0.1000 DL	0.1000	
10	13.640 ENDOSYL	-0.1000	-0.1000	226151.00	2.2615E-06	-0.0001 DL	-0.1000	
11	14.181 ENDOSYL ACID	-0.1000	-0.1000	110309.00	1.1030E-06	0.1000 DL	-0.1000	
12	14.371 DDT	1.0000	1.0000	10000.00	1.0000E-06	0.1000 DL	-0.1000	
13	17.430 PCB	0.1000	0.1000	18516.00	1.8516E-06	-0.1000 DL	0.1000	
Total Area: 41,0370 18703641.00 1.0000								

Varian 3400 Capillary Column GC: DB-1701 30M x 0.53mm I.D., Detector: 10000,
Injector: 250°C

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Sample Date : 10/17/2011 10:13:19

File Name : C:\Users\104143\OneDrive\HPLC\1017\1017.DAT

Start Time : 0.00 min

End Time : 30.00 min

Scale Factor : 0.1

Plot Offset : -100 mV

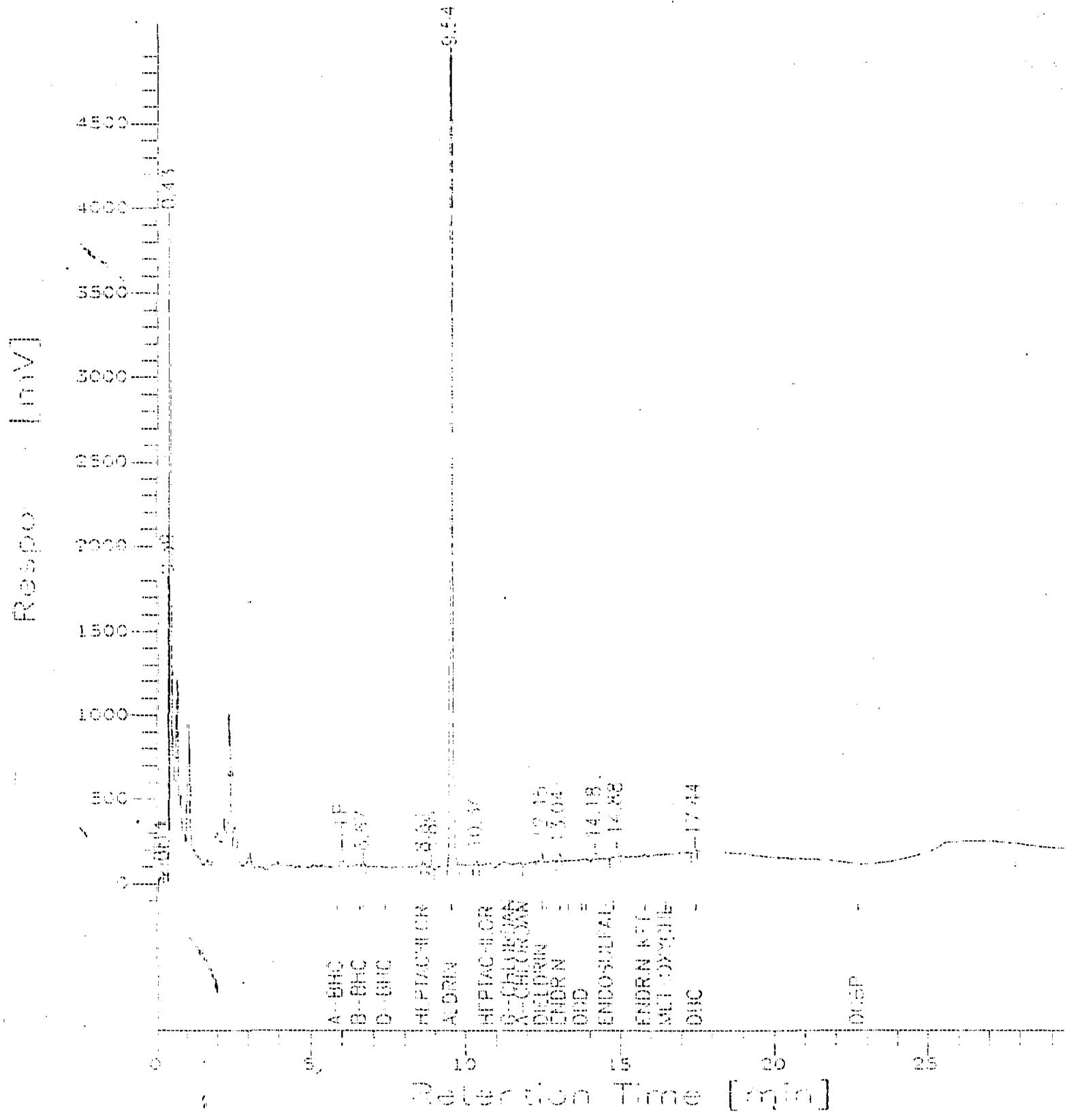
Sample ID : 1017

Date : 10/17/2011 10:13:19

Low Point : -100.00 mV

High Point : 4500.00 mV

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Software Version: 3.1 42425
 Date: 12/24/01 10:46
 Sample Name : 8860-3 XS Acid Cl
 Data File : d:\27301\data\1224017.daw Date: 12/24/01 10:46
 Sequence File: D:\27301\DATA\1224017.seq Cycles: 7 Channel: 1-A
 Instrument : VARIAN 3400 Pack/Width: Q/C Operator: 37
 Sample Amount : 1.0000 Dilution Factor : 1.00

Princeton Testing Laboratory Analysis Report

PESTICIDES AND POISNS BY CAPILLARY CHROMATOGRAPHY

Peak Ret Time 4 [min]	Component Name	Ref. Amount Concentration [ug/L]	Area	Height	St Dev	Rt	Reference	Cal. (Molar)	Range Factor
1 0.128		-1.3454	10702.41	1771.03	33	-----		-0.3011	
2 4.251		17.5767	17.5767	4081.00	1.044	31	-----	1.4731	
3 6.364	6-BHC	29.1764	29.1764	10154337.00	1.008	42 -0.1281	BEC	1.3931	
4 6.672	6-BHC	5.3479	5.3479	683207.00	1.008	39 -0.1281	BEC	0.1021e3	
5 6.787		1.3899	1.3899	315103.75	100308.01	39 -----		0.1412e3	
6 6.971		5.3279	5.3279	680542.50	367451.01	39 -----		1.0312e3	
7 7.103		5.3946	5.3946	4812527.00	173148.05	39 -----		1.1112e3	
8 7.475	6-BHC	35.9763	35.9763	6674738.00	1.008	31 1.4073	BEC	1.4073	
9 7.673		2.4878	2.4878	128421.00	282411.04	31 -----		1.1721e3	
10 7.746		0.6954	0.6954	1093072.00	238701.17	31 -----		1.3747e3	
11 7.893		0.6611	0.6611	1502197.00	1.528	31 -----		1.1631e3	
12 8.138		1.3163	1.3163	3632338.00	503125.10	31 -----		1.4817e3	
13 8.158		0.6597	0.6597	1727471.10	350495.08	31 -----		1.1513e3	
14 8.406		0.9381	0.9381	1617401.00	212273.00	31 -----		1.1070e3	
15 8.516	HEPTACHLOR	0.5670	0.5670	139550.00	198650.01	31 -0.1631	BEC	1.1631e3	
16 8.716		-0.2764	-0.2764	435229.75	146323.24	31 -----		-0.1631	
17 8.931		2.4522	2.4522	5556493.50	660162.05	31 -----		1.1420e3	
18 9.214		1.9071	1.9071	2022365.75	308493.41	31 -----		1.1631e3	
19 9.417		4.1164	4.1164	4794267.00	764361.01	31 -----		1.1420e3	
20 9.526	ALDRIN	1.1591	1.1591	2336203.50	347223.15	31 -0.1512	BEC	1.1512e3	
21 9.695		1.3753	1.3753	1790511.50	251950.04	31 -----		1.1631e3	
22 9.910		10.1369	10.1369	11992954.00	1.018	31 -----		1.1812	
23 10.140		5.3521	5.3521	5237570.50	1.028	31 -----		1.1631e3	
24 10.293		4.0325	4.0325	3560363.00	458463.01	31 -----		5.000e3	
25 10.505		13.9111	13.9111	13727175.00	1.008	31 -----		1.1631	
26 10.770	HEPTACHLOR EPO	0.5616	0.5616	8653821.00	1.138	31 0.5103	BEC	1.1631e3	
27 11.391		1.1925	1.1925	1429447.50	304957.06	31 -----		1.1631e3	
28 11.462		23.6346	23.6346	24167326.00	1.178	31 -----		1.4495	
29 11.570	6-CHLOROHE	3.3033	3.3033	1427431.10	278519.04	31 -0.1281	BEC	1.1280e3	
30 11.580		4.2742	4.2742	2293627.50	381701.47	31 -----		1.3730e3	
31 11.644		2.1567	2.1567	1759436.10	292873.49	31 -----		0.1532e3	
32 11.806	ENDO 1	0.9705	0.9705	535801.25	106882.57	31 -----		0.1512e3	
33 11.972	6-CHLOROHE	5.0655	5.0655	380266.19	130140.04	31 1.0510	KETOMONCHLOR	1.2437e3	
34 12.090		4.3365	4.3365	344121.72	52142.27	31 -----		19354.2912	
35 12.280		5.9838	5.9838	4834981.50	500035.31	31 -----		1.4112e3	
36 12.465	DIELDRIN	14.2284	14.2284	2798173.50	367868.47	31 -0.0413	BEC	0.6175e3	
37 12.761	ENDO 2	11.3552	11.3552	7953934.50	227717.13	31 1.5845	BEC	1.2134	
38 13.016	ENDRIN	11.5275	11.5275	9262067.00	566334.05	31 -0.1853	BEC	1.3047e3	
39 13.274	ENDO 2	-1.6745	-1.6745	1182938.63	233455.52	31 -0.2432	BEC	-0.2432	

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40	13.528	13.7833	11.7573	6503.64,15 3e4e13,25 W -----	4.112
41	13.723 900 "	21.1999	11.1999	2918711.00 412282.66 W -0.1413 0.00	4.0310e5
	13.734 ENCRIM ALDE	8.0354	6.0354	3591265.00 354353.47 W 0.4326 0.01	3.5107e1
	14.113	5.1323	5.1323	3026992.75 475013.53 W -----	5.3872e5
44	14.427 ENODISULFATE	17.6473	17.6473	17415628.00 1.6368 W -0.1207 0.01	4.6112
45	14.713 031	15.2114	15.2114	4771283.00 1.1268 W -0.4013 0.01	4.0310
46	14.752	21.3594	21.3594	1113227.00 534677.44 W -----	5.0301
47	15.193	14.4072	14.4072	7315011.00 674269.44 W -----	5.0300
48	15.340	3.4247	3.4247	1699411.75 196399.94 W -----	6.7023e5
49	15.495	0.1138	0.1138	535661.52 40154.61 W -----	2.4491e1
50	15.627	0.7395	0.7395	707136.00 103820.05 W -----	8.8433e1
51	15.713	3.2352	3.2352	2029361.00 303618.57 W -----	4.4311e1
52	15.870 ENDRATH REZONE	0.5324	0.5324	585237.07 117735.60 W 0.1544 0.01	4.1124e1
53	15.931	9.1071	9.1071	1924261.00 179471.61 W -----	2.8311e1
54	16.511	0.5748	0.5748	561614.00 111575.44 W -----	4.1124e1
55	16.414	0.6460	0.6460	351651.00 351651.00 W -----	1.0310e1
56	16.574 ENTRAPROCHLOR	21.5173	21.5173	612645.50 464113.15 W 1.4211 0.01	2.1124e1
57	16.732	0.0361	0.0361	1774454.00 158101.80 W -----	1.4310e1
58	17.631	1.8411	1.8411	820375.00 203870.00 W -----	4.1124e1
59	17.237	0.3193	0.3193	231401.00 60361.00 W -----	4.1124e1
60	17.411	0.2284	0.2284	190495.00 140495.00 W -----	4.1124e1
61	17.535 030	0.6571	0.6571	2946101.00 474601.00 W 0.3000 0.01	4.1124e1
62	17.536	1.0708	1.0708	921401.00 142101.00 W -----	4.1124e1
63	17.107	0.0987	0.0987	178754.00 158101.80 W -----	4.1124e1
64	17.123	0.3384	0.3384	159701.00 65757.00 W -----	4.1124e1
65	17.421	0.1229	0.1229	54801.00 54801.00 W -----	4.1124e1
66	17.468	0.5153	0.5153	372650.00 17571.00 W -----	4.1124e1
67	17.501	0.8061	0.8061	113604.00 80445.00 W -----	4.1124e1
68	17.523	0.5227	0.5227	2097461.00 146501.00 W -----	4.1124e1
69	17.524	0.0617	0.0617	178011.00 13479.00 W -----	4.1124e1
70	17.547	0.2453	0.2453	77321.00 10541.00 W -----	4.1124e1
71	17.551	1.6077	1.6077	449301.00 17357.00 W -----	4.1124e1
72	17.561	0.7819	0.7819	437682.00 17354.00 W -----	4.1124e1
73	17.581	0.4079	0.4079	181879.00 19689.00 W -----	4.1124e1
74	20.157	15.7726	15.7726	624631.50 523461.50 W -----	4.1124e1
75	20.401	-0.5725	-0.5725	512264.50 105629.00 W -----	4.1124e1
76	23.151 03P	-2.2302	-2.2302	140721.00 21221.00 W 1.9310 0.01	4.1124e1
77	23.703	0.9668	0.9668	773577.00 63461.00 W -----	4.1124e1
78	24.056	-2.1229	-2.1229	140413.50 16742.00 W -----	4.1124e1
79	24.058	-0.5861	-0.5861	75470.00 4174.00 W -----	4.1124e1
80	25.227	-0.1332	-0.1332	10303.00 16101.00 W -----	4.1124e1
81	27.032	2.0318	2.0318	310159.00 158731.00 W -----	4.1124e1

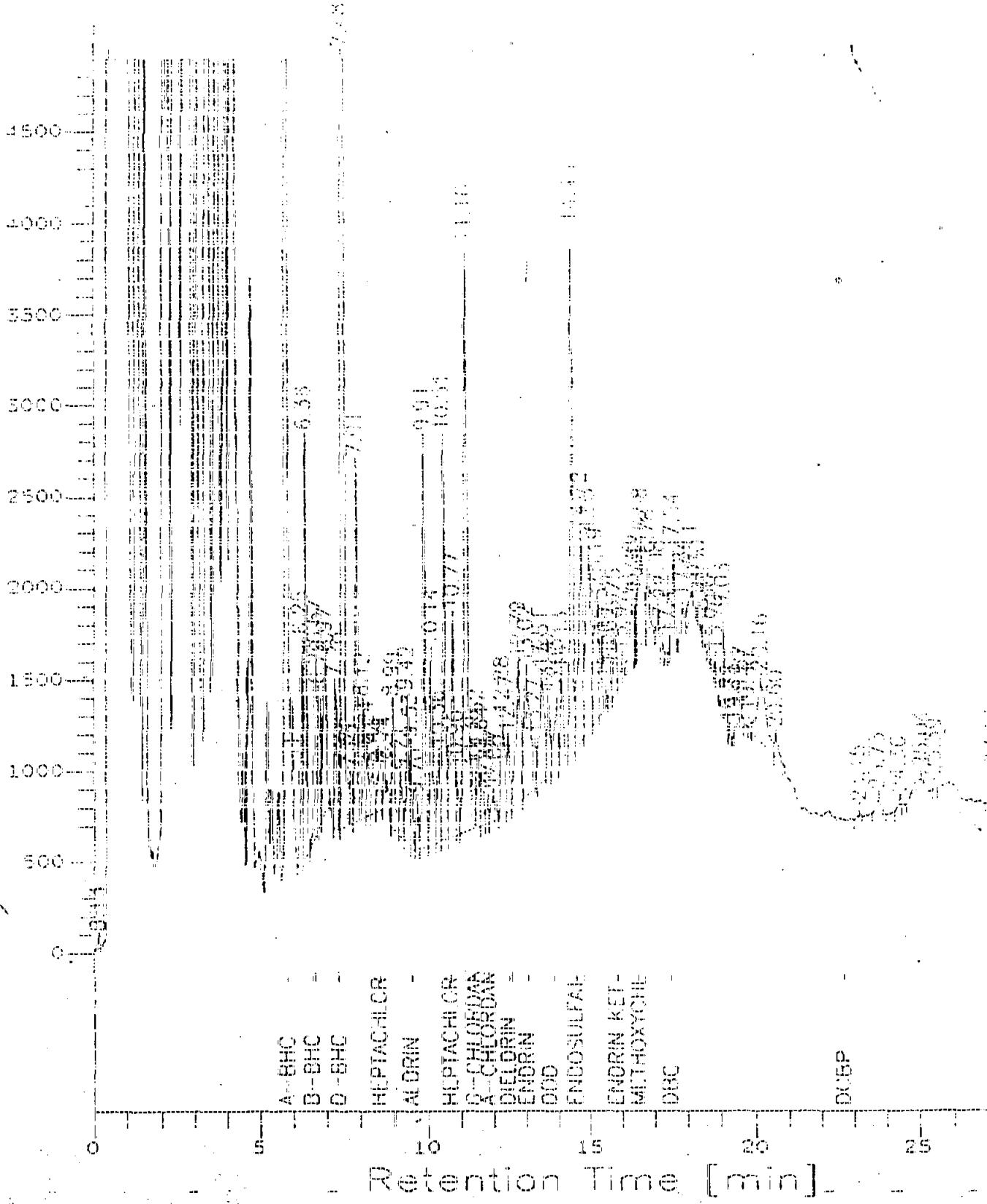
461.073 461.073 1.000 1.000

Marian 3400 Capillary Column 60:1 0.01 COM x 0.01mm I.D. Detector: 250C
Injector: 250 C

AR100641

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Response [mV]



Retention Time [min]

BR100642

00009

Software Version: 5.1 - DALS

Date: 11/21/91 File: 51031

Sample Name : 1001 SCUGAL

Data File : d:\ptl\001\data\1001.daw Date: 11/21/91 8:21

Source File: D:\PTL\DATA\1001.daw File: 51031.DAT

Instrument : VARIAN 3400 - Pack. Date: 9-6-91 Operator: A0

Sample Amount : 1.0000 Dilution Factor : 1.00

Princeton Testing Laboratory Analysis Report

PESTICIDES AND PCB'S BY CAPILLARY CHROMATOGRAPHY

Peak Ret. Time T [min]	Concentr. ppm	Raw Average Concentration [ug/l]	Area [uV-sec]	Height [uV]	Delta H [uV]	Reference Concentr.	Cal. Intercept	Cal. Slope
1	0.100	0.0000	340000.00	30610.72	00	-----	-----	1.0000e-0
2	0.200	0.0000	350451.03	6.525.40	00	-----	-----	1.0000e-0
3	0.400	1.0000	1071571.05	307143.01	00	-----	-----	1.0000e-0
4	0.600	0.1178	177364.07	145501.42	07	-----	-----	1.0000e-0
5	0.800	0.0000	2903850.00	330253.00	00	-----	-----	1.0000e-0
6	1.000	0.0000	507744.00	54791.06	02	-----	-----	1.0000e-0
7	1.100	0.4200	426532.00	42295.50	00	-----	-----	1.0000e-0
8	1.300	0.0000	810944.00	43475.00	00	-----	-----	1.0000e-0
9	14.000	0.0000	865864.00	103671.40	00	-----	-----	1.0000e-0
10	14.710	0.0000	811161.00	107479.72	00	-----	-----	1.0000e-0
		0.0000	3286101.00	1.4946				

Variian 3400 Capillary Column GC: DB-5CB 30M x 0.25mm I.D. Detector: ECD.

Injector: 250°C

AR100643

00040

Pesticide-POC Chromatogram

Sample name : 1001.D0001

File name : 1001.D0001.P0001.DAT

Start Time : 1.00 min End Time : 10.00 min

Scale Factor : 1 Plot Offset : 100.00

Sample ID : 1001

File date : 01/01/01

File time : 10:00:00

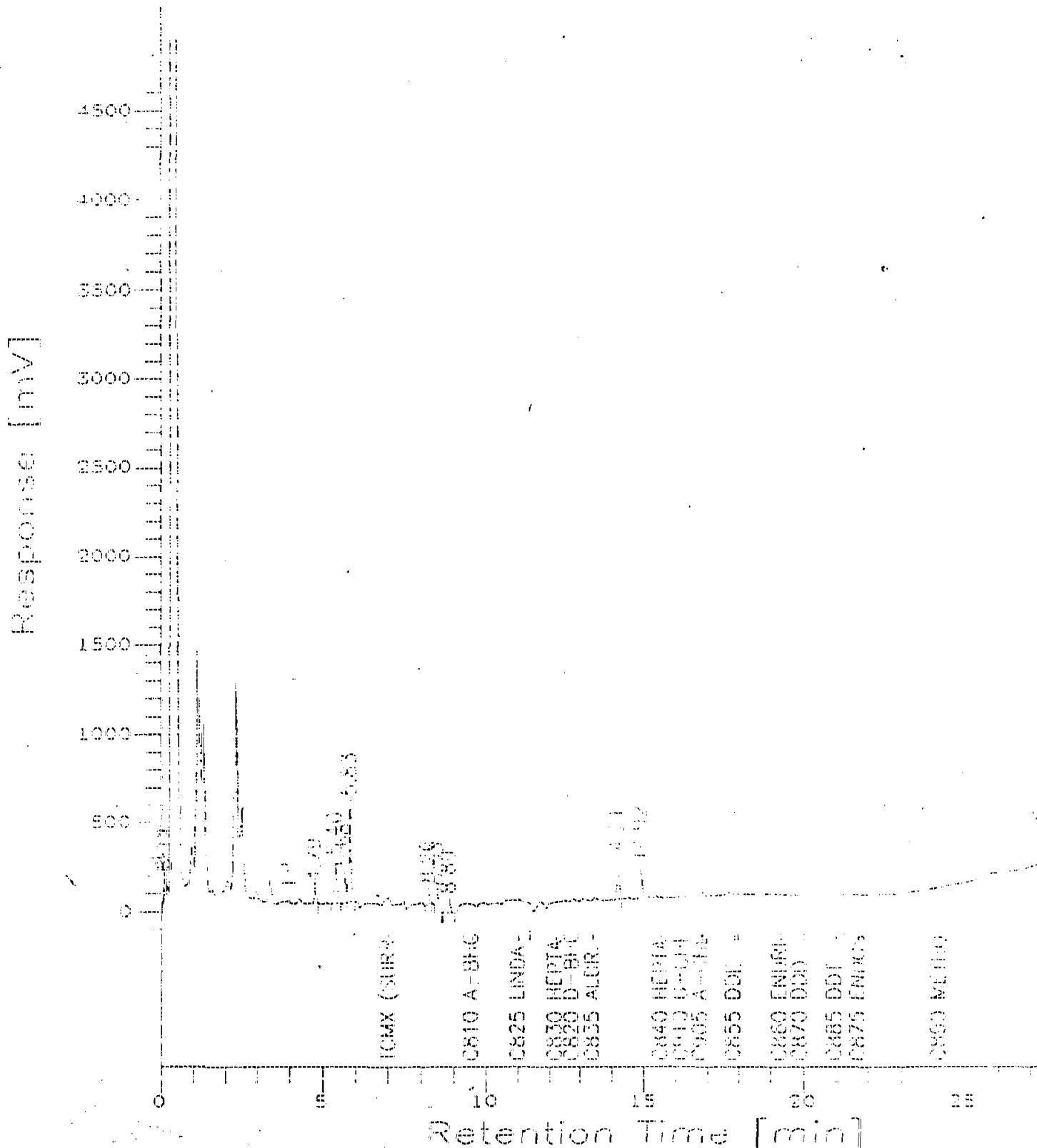
Plot offset : 100.00

Sample ID : 1001

File date : 01/01/01

File time : 10:00:00

Plot offset : 100.00



AR100644

00011

Sample ID: 100645-A
Sample Name: 100645-A

Matrix: Water
Sample File: 100645-A.DAT
Reference File: 100645-A.DAT
Instrument: Varian 3400 GC
Date: 10/18/2012
Time: 10:34:45
Run Time: 00:00:00
Chromatogram Type: Total Ion Chromatogram
Integration Factor: 1.000000

Princeton Testing Laboratory Analytical Report

100645-A.DAT 10/18/2012 10:34:45 TIC.DAT 10/18/2012

Retention Time min	Component #	Estimated Concentration ppm	Area	Relative Retention Time	Library ID	Reference	Calibration Factor
1.011		0.000	0.0000	0.0000	000000000000	-----	1.00000
1.136		1.118	1.1180	0.1180	000000000000	-----	1.00000
1.140		1.188	1.1880	0.1880	000000000000	-----	1.00000
1.157		1.117	1.1170	0.1170	000000000000	-----	1.00000
1.224		4.224	4.2240	0.2240	000000000000	-----	1.00000
1.317 (100% Peak)		1.317	1.3170	0.3170	000000000000	-----	1.00000
		0.411	0.4110	0.4110	000000000000	-----	1.00000
		0.381	0.3810	0.3810	000000000000	-----	1.00000
		0.374	0.3740	0.3740	000000000000	-----	1.00000
		0.376	0.3760	0.3760	000000000000	-----	1.00000
		0.733	0.7330	0.7330	000000000000	-----	1.00000
		0.813	0.8130	0.8130	000000000000	-----	1.00000
		0.814	0.8140	0.8140	000000000000	-----	1.00000
		0.829	0.8290	0.8290	000000000000	-----	1.00000
		0.840 (100% Peak)	0.840	0.840	000000000000	-----	1.00000
		0.853	0.8530	0.8530	000000000000	-----	1.00000
		0.941	0.9410	0.9410	000000000000	-----	1.00000
		10.463 (100% Peak)	10.463	10.463	000000000000	-----	1.00000
		10.474	10.4740	10.4740	000000000000	-----	1.00000
		10.535	10.5350	10.5350	000000000000	-----	1.00000
		10.539	10.5390	10.5390	000000000000	-----	1.00000
		10.465	10.4650	10.4650	000000000000	-----	1.00000
		10.466 (100% Peak)	10.466	10.466	000000000000	-----	1.00000
		10.477	10.4770	10.4770	000000000000	-----	1.00000
		10.481	10.4810	10.4810	000000000000	-----	1.00000
		10.481	10.4810	10.4810	000000000000	-----	1.00000
		10.481	10.4810	10.4810	000000000000	-----	1.00000
		10.481 (100% Peak)	10.481	10.481	000000000000	-----	1.00000

Total Run Time: 00:00:00.000000

Instrument: Varian 3400 Capillary Column: G1: 08-008 30M X 0.32MM ID ID. Detector: 3000.

Injector: 300 C

18680
AR100645

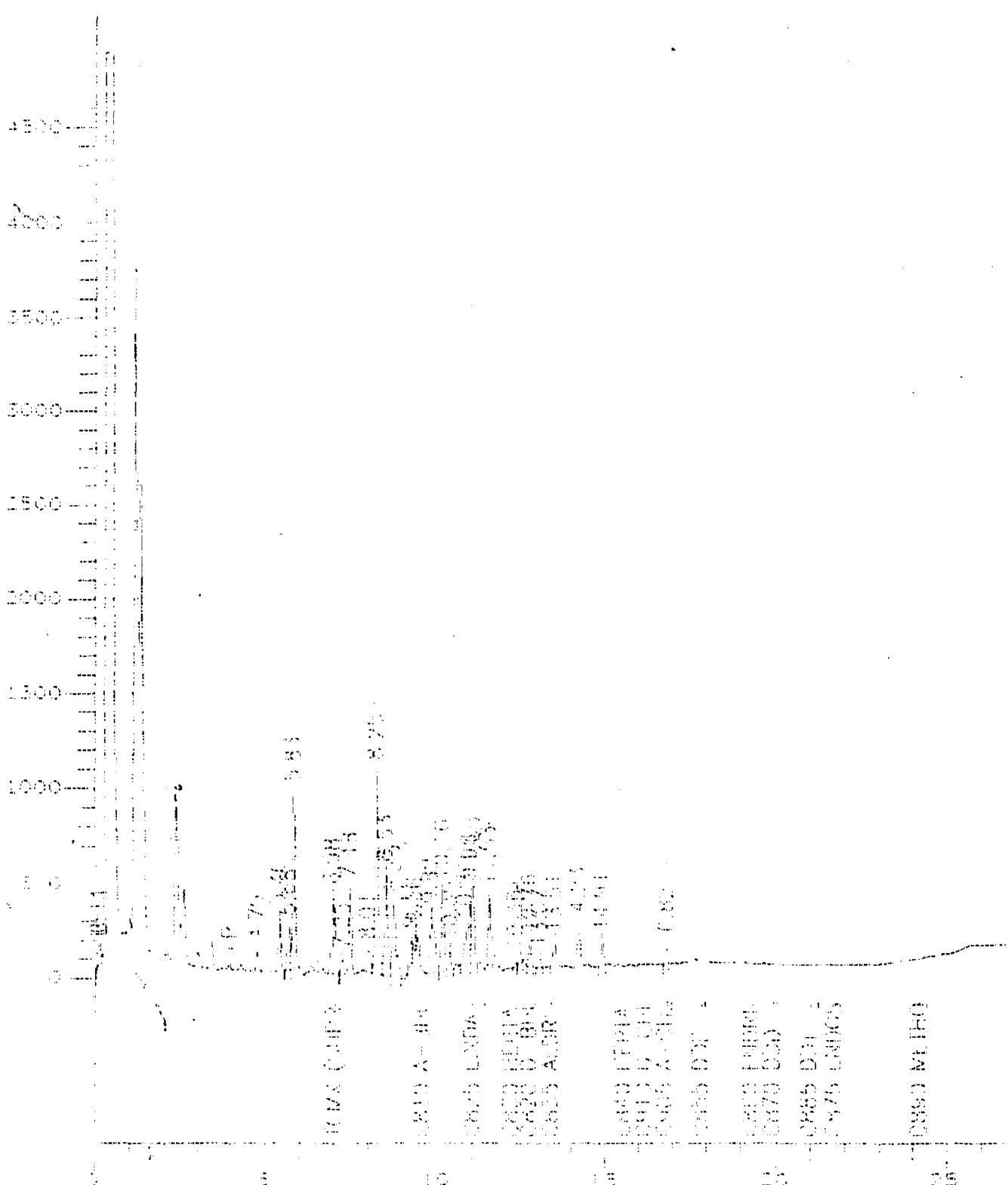
* 00012

Waters HPLC Data Processing Report

Sample Name : 1032 500001
Filename : d:\data\detected\1032\1032.WDX
Start Time : 0.00 min End Time : 10.00 min
Scale Factor: 0 Plot Offset: -100.00 mV

Sample #: 21 Date : 11/21/01 6:02
Low Point : -100.00 mV High Point : 1800.00 mV
Plot Scale: 5000 mV

Page 1 of 1



Waters HPLC Data Processing Report

AR100646

0001

Software Version: 3.1 RELEASE

Date: 11/21/91 Time:

Sample Name : 1012 3000ML

Data File : C:\PTCO\DATA\1121020.D Date: 11/21/91 ext: D

Sequence File: C:\PTCO\DATA\1121020.seq File: 02 Infrared: 04

Instrument : VARIAN 3400 Rack/Visit: C,1 Operator: 4K

Sample Amount : 1.0000 Collection Factor : 1.00

Princeton Testing Laboratory Analysis Report

PESTICIDES AND PCB'S BY CAPILLARY CHROMATOGRAPHY

Peak Ret. Time T (min)	Component Name	Avg Abstat Concentration (ppm) (ppb)	Area (mV-sec)	Relative to Delta T (%)	Reference Compound	Rel. Calibrated Abst. Factor	
1	0.113	0.0147	0.1000	224701.34	100.00	00	1.0000e1
2	4.803	0.0113	0.0000	111807.01	41.91	00	1.0000e0
3	5.414	0.0105	0.0000	381479.02	21.43	00	1.0000e0
4	5.538	0.0010	0.0000	323608.00	1.00	00	1.0000e0
5	5.839	0.0006	0.0000	233730.00	4.00	00	1.0000e0
6	6.051	0.1954	0.0000	103339.05	3.00	00	1.0000e1
7	6.979 1012 3000ML	0.0000	0.0000	324056.06	49.10	1012 3000	1.0000e1
8	7.158	0.0017	0.0000	201061.02	6.00	00	1.0000e0
9	7.481	0.0452	0.0000	2245161.03	46.74	00	-1.0000e0
10	8.002	0.0779	0.0000	375161.03	8.00	00	1.0000e0
11	8.247	0.4110	0.0000	6413781.00	1.00	00	1.0000e0
12	8.503	0.9751	0.0000	2970031.01	6.00	00	1.0000e0
13	8.738	0.6504	0.0000	2410681.01	47.00	00	1.0000e0
14	9.235	0.1661	0.0000	494181.01	15.00	00	1.0000e0
15	9.356	0.0841	0.0000	1024123.00	20.00	00	1.0000e0
16	9.453 0312 4-BHC	1.2261	0.0000	1011214.06	100.00	1012 3000	1.0000e0
17	9.811	0.0000	0.0000	10330119.00	20.00	00	1.0000e0
18	10.147	0.1770	0.0000	2472956.00	4.00	00	1.0000e0
19	10.379	0.3772	0.0000	361160.04	8.00	00	1.0000e0
20	10.529	0.1360	0.0000	585977.00	27.00	00	1.0000e0
21	10.943 0312 4-BHC	1.5527	0.0000	755820.04	14.00	1012 3000	1.0000e0
22	11.045	0.5762	0.0000	575030.04	18.00	00	1.0000e0
23	11.512	0.1650	0.0000	162326.00	3.00	00	1.0000e0
		15.0349	0.0000	31480301.00	1.00		

Varian 3400 Capillary Column GC: DB-500 30m x 0.25mm I.D. Detector: 3000C

Injector: 250 C

AR100647

0001

Pesticide A/C/E Chromatogram

Sample Name : 1015 300601

File Name : 10150601.DAT\1015.DAT

Start Time : 0.00 min End Time : 20.00 min

Scale Factor : 0 Peak Offset : 0.00 mV

Sample #: 10

Date : 11/11/01 10:10

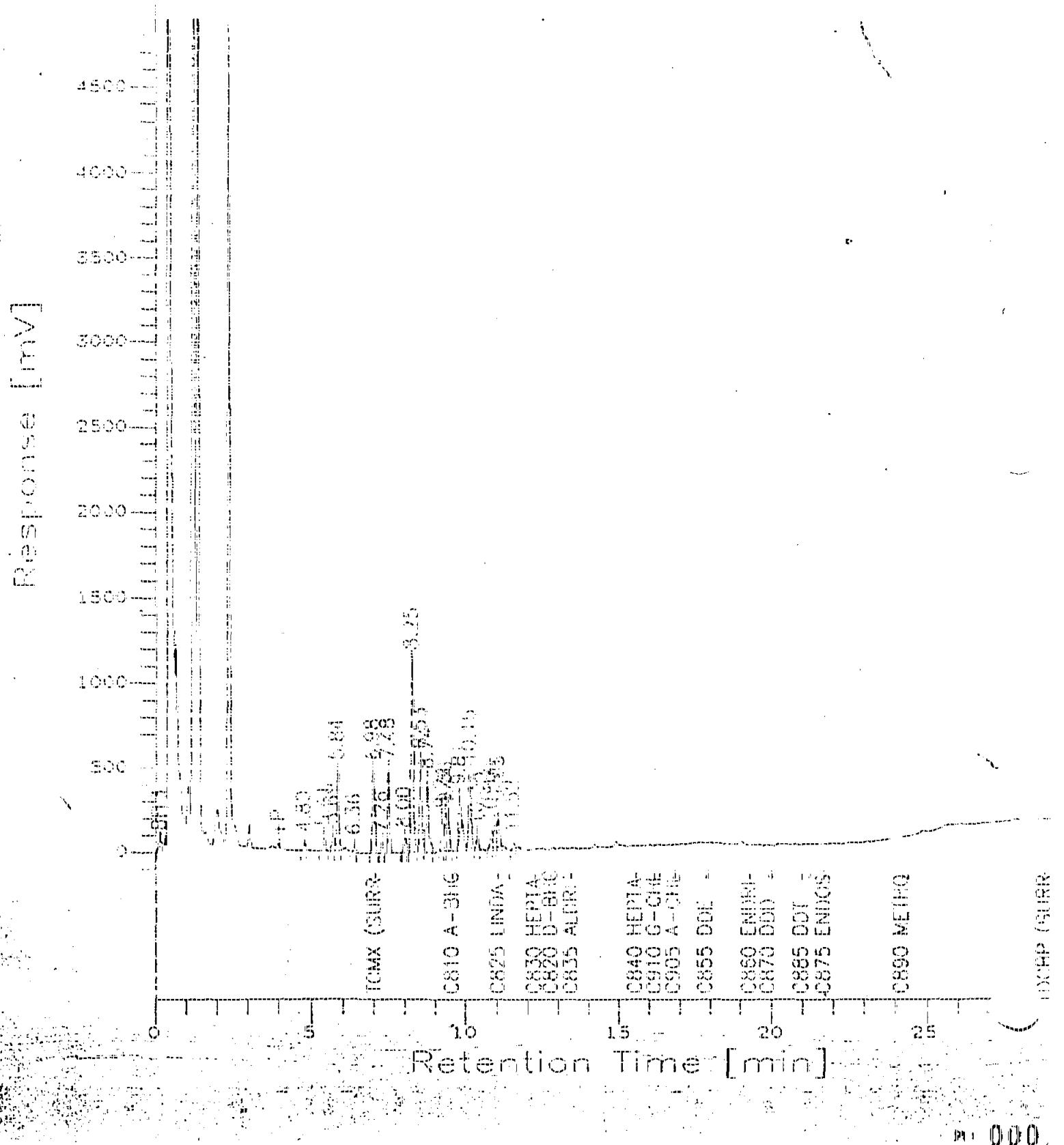
End Point : 20.00 min

Plot Scale : 400.00 mV

Page 1 of 1

Sign. Scale : 400.00 mV

Plot Scale: 1000 mV



Software Version: 3.1.0.402

Date: 11/21/01 7:56

Sample Name: 10241-400G-1

Data File: d:\ptc\data\10241-400G-1.DAT Date: 11/21/01 7:56

Source File: d:\ptc\data\10241-400G-1.DAT Scan: 10241-400G-1

Instrument: VARIAN 3400 Rack Wklt: G/C Detector: FID

Sample Amount: 1.0000 Dilution Factor: 1.00

Princeton Testing Laboratory Analysis Report

PESTICIDES AND PCB'S IN CAPILLARY CHROMATOGRAPHY

Peak Ret. Time t (min)	Component Name	Raw Absorbance Concentration (ppm)	Area	Height	R. Delta H	Reference	Call. Calibration Factor
1	0.108	0.0001	0.000000.00	0.000000.00	-----		1.0000e0
2	0.109	0.0002	0.000000.00	0.000000.00	-----		1.0000e0
3	0.408	0.0010	0.000000.00	0.000000.00	-----		1.0000e0
4	0.409	0.0010	0.000000.00	0.000000.00	-----		1.0000e0
5	0.444	0.0009	0.000000.00	0.000000.00	-----		1.0000e0
6	0.565	0.0009	0.000000.00	0.000000.00	-----		1.0000e0
7	0.878 Toluene	0.0010	0.000000.00	0.000000.00	0.000000.00	Toluene	1.0000e0
8	1.193	0.0027	0.000000.00	0.000000.00	-----		1.0000e0
9	1.433	0.0009	0.000000.00	0.000000.00	-----		1.0000e0
10	1.875	0.0001	0.000000.00	0.000000.00	-----		1.0000e0
11	2.101	0.0011	0.000000.00	0.000000.00	-----		1.0000e0
12	2.179	0.0020	0.000000.00	0.000000.00	-----		1.0000e0
13	2.342	0.0006	0.000000.00	0.000000.00	-----		1.0000e0
14	2.753	0.0006	0.000000.00	0.000000.00	-----		1.0000e0
15	3.003	0.0006	0.000000.00	0.000000.00	-----		1.0000e0
16	3.441 Dieldrin	0.0004	0.000000.00	0.000000.00	0.000000.00	Dieldrin	1.0000e0
17	3.815	0.0006	0.000000.00	0.000000.00	-----		1.0000e0
18	10.132	0.0005	0.000000.00	0.000000.00	-----		1.0000e0
19	10.381	0.0005	0.000000.00	0.000000.00	-----		1.0000e0
20	10.656	0.0002	0.000000.00	0.000000.00	-----		1.0000e0
21	10.952 ODEE Lindane	0.0006	0.000000.00	0.000000.00	0.000000.00	Lindane	1.0000e0
22	11.046	0.0004	0.000000.00	0.000000.00	-----		1.0000e0
23	11.504	0.0006	0.000000.00	0.000000.00	-----		1.0000e0
24	11.882	0.0004	0.000000.00	0.000000.00	-----		1.0000e0
25	11.935	0.0006	0.000000.00	0.000000.00	-----		1.0000e0
26	11.463	0.0001	0.000000.00	0.000000.00	-----		1.0000e0
27	11.754 Dieldrin	0.0001	0.000000.00	0.000000.00	0.000000.00	Dieldrin	1.0000e0
28	17.052	0.0003	0.000000.00	0.000000.00	-----		1.0000e0
29	17.511 ODEE Lindane	0.0003	0.000000.00	0.000000.00	0.000000.00	Lindane	1.0000e0
30	14.249	0.0007	0.000000.00	0.000000.00	-----		1.0000e0
31	14.807	0.0005	0.000000.00	0.000000.00	-----		1.0000e0

49.9339 13.4620 0.019424.00 1.95e4

Instrument: VARIAN 3400 Capillary Column GC: DB-50B 30M x 0.53mm I.D. Detector: FID
Injector: 250 °C

ARI 00649

00016

Pesticides/PCB Chromatogram

Sample Name : 1201-400001

FileName : d:\21000\data\3\1121013.rev

Start Time : 0.00 min

End Time : 30.00 min

Scale Factor : 6

Plot Offset: -100.00

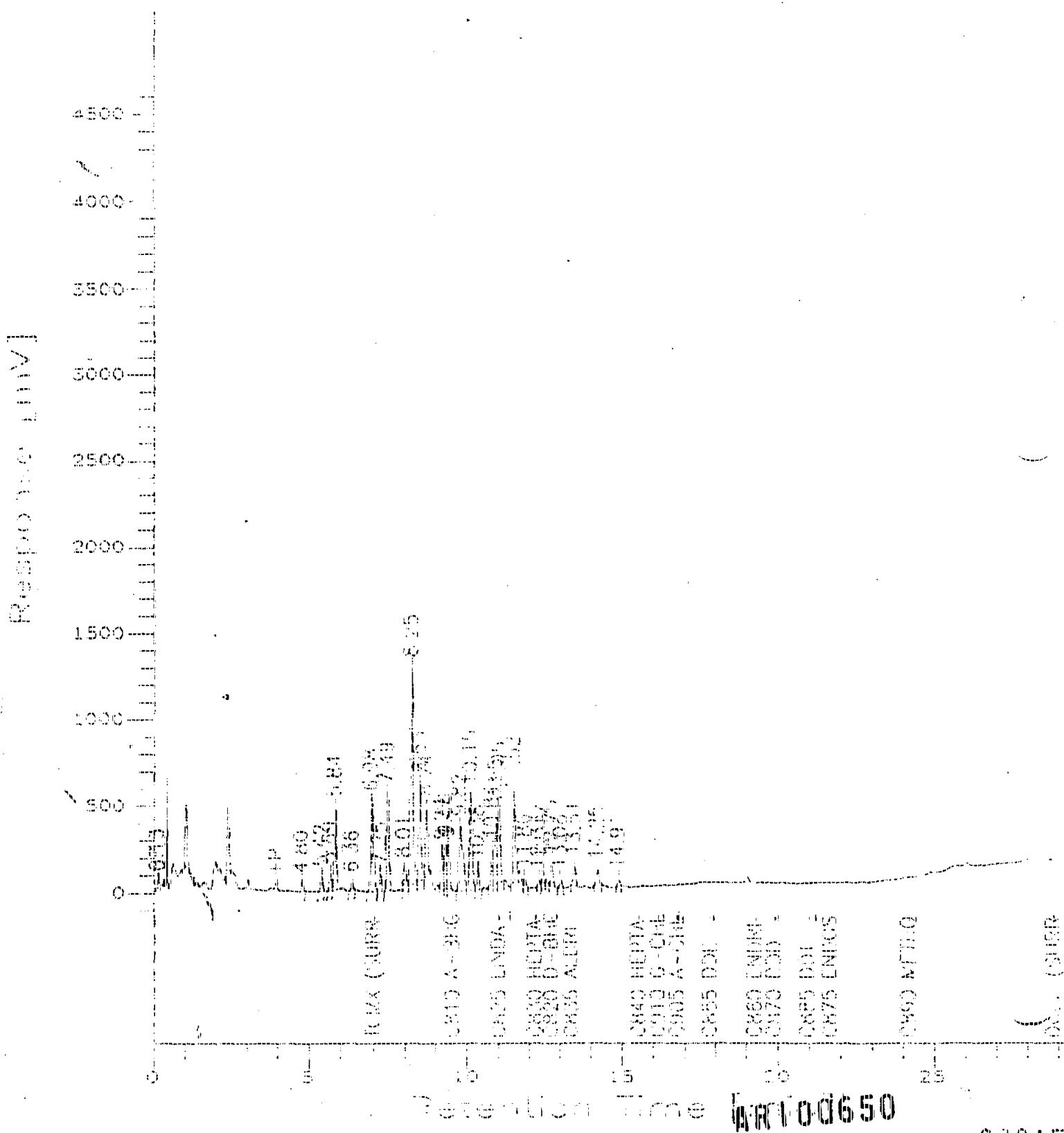
Sample ID: 21

Date : 11/21/01 11:37

Low Point : -100.00 mV

High Point : 4000.00 mV

Page 1 of 1



0.0000 MPPDQ

00017

Software version: 3.1 <2A23>

Date: 11/21/91 7:41

Sample Name : 1248 SOCUG/L

Data File : d:\3700\data\1121024.raw Date: 11/21/91 7:41

Sequence File: C:\3700\DATA\11214.seq File: 24 Channels: 4

Instrument : VARIAN 3400 Rack/Vials: 0/0 Operator: AK

Sample Amount : 1.0000 Dilution Factor : 1.00

Princeton Testing Laboratory Analysis Report

PESTICIDES AND POS'S BY CAPILLARY CHROMATOGRAPHY

Peak Ret Time # (min)	Component Name	Raw Present Concentration [ug/l]	Net [ug/l]	Area [mV-sec]	Height: H1 Delta F ^a [mV]	Reference Detector	Cal. Calibrated Range Factor
1	0.115	0.3416	0.0000	341654.06	62954.07	H6 -----	1.0000e
2	6.550 DDE (DDE)	0.5395	0.0005	1091881.03	173861.84	H6 -1.0000	1.0000e
3	7.456	0.7606	0.0000	1075361.61	161865.97	H6 -----	1.0000e
4	8.118	1.5593	0.0000	5557001.75	814614.69	H6 -----	1.0000e
5	8.557	1.5887	0.0000	1050711.75	211774.41	H6 -----	1.0000e
6	8.750	0.7602	0.0000	760204.06	131151.17	H6 -----	1.0000e
7	8.856	0.7782	0.0000	778075.06	112044.31	H6 -----	1.0000e
8	9.223	0.4504	0.0000	336937.44	114230.28	H6 -----	1.0000e
9	9.343 ODE A-EMC	0.7415	0.7313	104035.44	10380.77	H6 -1.0000	1.0000e
10	9.801	0.4594	0.0000	1613066.33	340101.50	H6 -----	1.0000e
11	10.144	0.4493	0.0000	3465913.03	592301.23	H6 -----	1.0000e
12	10.372	0.6711	0.0000	401187.75	151842.21	H6 -----	1.0000e
13	10.623	1.3575	0.0000	1077451.75	110740.56	H6 -----	1.0000e
14	10.949 1035 LINDBE	0.8182	0.0002	1074818.06	1070111.01	H6 -0.1000	1.0000e
15	11.037	0.4673	0.0000	1647349.25	460186.48	H6 -----	1.0000e
16	11.155	0.6702	0.0000	1570861.05	379401.15	H6 -----	1.0000e
17	11.258	1.6102	0.0000	1612844.05	1070101.84	H6 -----	1.0000e
18	11.656	1.4761	1.3000	471181.81	37380.84	H6 -----	1.0000e
19	12.315	0.6018	0.2000	601747.09	105027.41	H6 -----	1.0000e
20	12.451	0.4455	0.0000	1440745.05	213610.50	H6 -----	1.0000e
21	12.576 ODE B-EMC	2.4707	2.4707	712045.45	142153.41	H6 -1.0000	1.0000e
22	12.741	1.1855	0.0000	2155513.10	373401.18	H6 -----	1.0000e
23	13.030 ODE ALDRIN	1.7648	1.7648	148221.25	113738.33	H6 -1.0000	1.0000e
24	17.495	1.2671	0.0000	265027.05	172054.18	H6 -----	1.0000e
25	18.235	1.2634	0.0000	2631571.50	63771.57	H6 -----	1.0000e
26	18.633	1.1412	0.0000	1141151.15	111611.50	H6 -----	1.0000e

40.1907 4.9501 41919716.00 7.0360

Varian 3400 Capillary Column GC: DB-608 30M X 0.53mm I.D. Detector: 3000.

Injector: 250 C

Pesticide/PCB Chromatogram

Sample Name : 1243 008671

File Name : 11\2300\adme3\1121024.rev

Start Time : 0.00 min End Time : 50.00 min

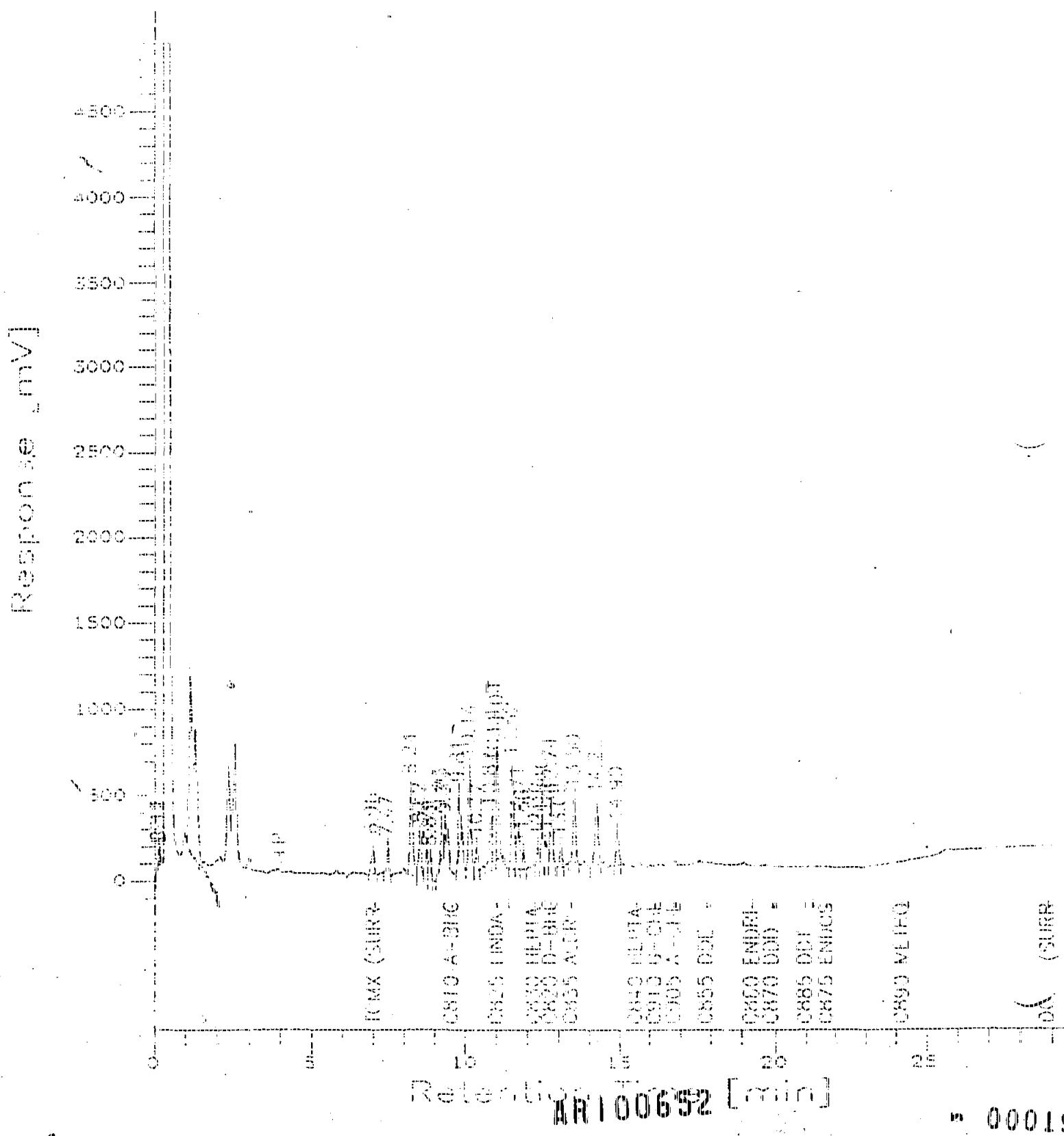
Scale Factor: 6 Plot Offset: -100.00 mV

Sample #: 21

Date : 11/21/91 8:11

Low Point : -100.00 mV High Point : 4900.00 mV

Plot Scales: 5000 mV



Software Version: 3.1 (2023)
Date: 11/21/91 9:55

Sample Name : CHLORDANE 1500g/L

Data File : c:\2700\data\1121027.han Date: 11-21-91 9:55
Sequence File: C:\2700\DATA\1121027.ser Cycle: 07 Channel: 2
Instrument : VARIAN 3400 Rack/Vial: 01 Operator: 4
Sample Amount : 1.0000 Dilution Factor : 1.00

Princeton Testing Laboratory Analysis Report

PESTICIDES AND PCB'S BY CAPILLARY CHROMATOGRAPHY

Peak Ret. Time # (min)	Component Name	Raw Amount Concentration (ug/L)	Raw Area (uV-sec)	Retention Time (min)	Delta RT (min)	Reference Component	Rel. Response Factor
1	0.132	0.0000	229817.40	00001.00	00	-	1.00000
2	0.155	0.0000	950531.00	00001.00	00	-	1.00000
3	0.190	0.0000	154514.00	00001.00	00	-	1.00000
4	0.237	0.0000	214269.00	00001.00	00	-	1.00000
5	0.245 TCM (SURR)	1.0000	190142.00	00001.00	00	TCM (SURR)	1.00000
6	0.256	0.0000	153776.00	00001.00	00	-	1.00000
7	0.263	0.0000	192391.00	00001.00	00	-	1.00000
8	0.292	0.0000	4464172.00	00001.00	00	-	1.00000
9	0.303	0.0000	5000197.00	00001.00	00	-	1.00000
10	0.344	0.0000	549305.00	00001.00	00	-	1.00000
11	0.355	0.0000	151104.00	00001.00	00	-	1.00000
12	0.366	0.0000	501073.00	00001.00	00	-	1.00000
13	0.378 TCM A-8PC	1.0000	1881706.00	00001.00	00	TCM (SURR)	1.00000
14	0.394	0.0000	390311.00	00001.00	00	-	1.00000
15	0.472	0.0000	215577.00	00001.00	00	-	1.00000
16	0.516	0.0000	3110570.00	00001.00	00	-	1.00000
17	10.970	0.0000	10137.00	00001.00	00	-	1.00000
18	11.139 CHLOR LINDANE	1.0000	487275.00	00001.00	00	TCM (SURR)	1.00000
19	11.153	0.0000	1423417.00	00001.00	00	-	1.00000
20	11.178	0.0000	1014514.00	00001.00	00	-	1.00000
21	11.632 CHL B-8PC	1.0000	1150280.00	00001.00	00	TCM (SURR)	1.00000
22	11.666	0.0000	9500397.00	00001.00	00	-	1.00000
23	12.009	0.0000	6471617.00	00001.00	00	-	1.00000
24	12.403 CHLO HEPTACHLOR	1.0000	712921.00	00001.00	00	TCM (SURR)	1.00000
25	12.513	0.0000	462176.00	00001.00	00	-	1.00000
26	12.542 CHLO D-8PC	1.0000	381681.00	00001.00	00	TCM (SURR)	1.00000
27	12.574 CHLO H-8PC	1.0000	1510464.00	00001.00	00	TCM (SURR)	1.00000
28	13.931	0.0000	2040861.00	00001.00	00	-	1.00000
29	14.212	0.0000	415512.00	00001.00	00	-	1.00000
		66.4950	19.7163	55519880.00	1.05e7		

Varian 3400 Capillary Column GC: DB-608 30M X 0.53mm I.D. Detector: 300C,
Injector: 250 C

AR100653

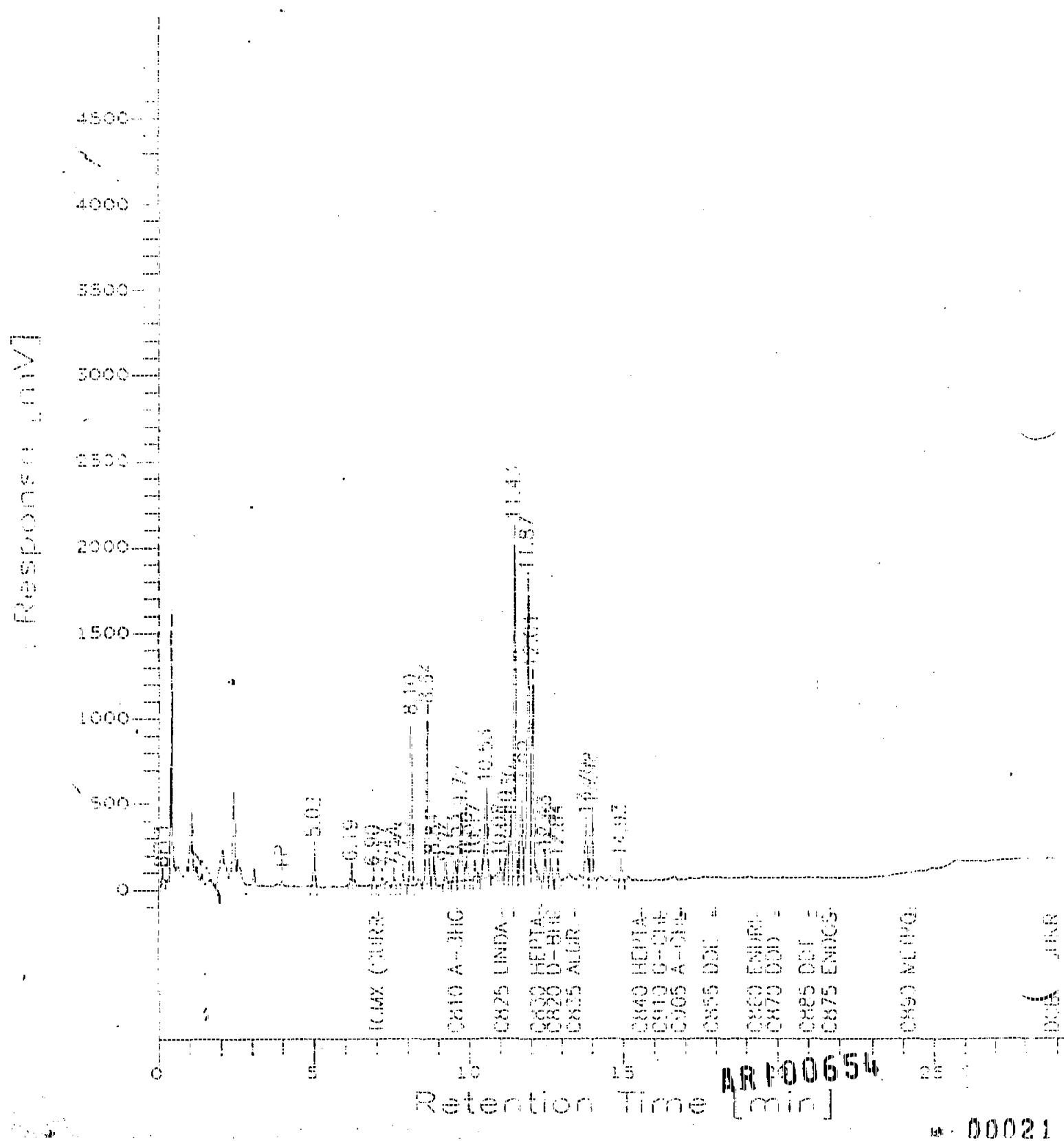
00029

Pesticide/PCB Chromatogram

Sample Name : CHLORGANE 15000/L
File Name : d:\2700\date\11121001.rws
Start Time : 0.00 min End Time : 50.00 min
Scale Factor : 0 Plot Offset : -100 mV

Sample ID: 27 Date: 11/21/91 9:56
Low Point : -100.00 mV High Point : 4500.00 mV
Plot Scale: 5000 mV

Page 1 of 1



Instrumental analysis was performed on a Varian 3300 GC/MS system.

Software Version 1.1, Date

Acquisition 12/10/1998

Sample Name 1121046

Data File 1121046.D, Date 12/10/98, Owner 1121046, Job 1121046

Sample File 1121046.D, Date 12/10/98, Owner 1121046, Job 1121046

Instrument Varian 3300, Reactor type C, Detector MS

Sample Amount 1.0000

Sample Type 1121046, Purity 1.0000

Sample Name 1121046, Purity 1.0000

Princeton Testing Laboratory Analysis Report

PESTICIDE AND POLY(1,4-PHENYLENE TEREPHTHALATE) ANALYSIS

Peak Retention t [min]	Chemical Name	Ref. Spectra Concentration [ug/L]	Area	Height at Delta H [uV]	Reference	Cell. Aliqnt	Form Factor
1	1,114	1.1764	80845.00	11351.61	-----	-----	4.000e5
2	0.108	34.7407	34.7407	34.7407	-----	-----	1.000
3	0.417 4-BEC	4.7401	4.7401	44935.00	11351.61	10000.000	-----
4	1.014	-0.1019	-0.1019	41036.00	11351.61	-----	4.000e5
5	0.104	-0.1019	-0.1019	10194.00	11351.61	-----	4.000e5
6	0.108 4-BENZYLIC	-0.1009	-0.1009	41036.00	11351.61	10000.000	-----
7	0.103	0.1043	0.1043	10194.00	11351.61	-----	4.000e5
8	0.103	0.1043	0.1043	10194.00	11351.61	-----	4.000e5
9	0.103 4-EPOXY	0.1034	0.1034	10200.00	11351.61	10000.000	-----
10	0.103	1.1497	1.1497	10177.00	11351.61	-----	4.000e5
11	10.112	1.1410	1.1410	16451.00	11351.61	-----	4.000e5
12	10.114	1.1392	1.1392	16511.00	11351.61	-----	4.000e5
13	10.107 4-PHENYL 4-EDO	1.1240	1.1240	41036.00	11351.61	10000.000	-----
14	10.108	1.1202	1.1202	10194.00	11351.61	-----	4.000e5
15	11.012	4.7391	4.7391	41036.00	11351.61	-----	4.000e5
16	11.010	1.1380	1.1380	10194.00	11351.61	-----	4.000e5
17	11.562 4-CHLOROCINE	1.0714	1.0714	10194.00	11351.61	10000.000	-----
18	11.720 ENOD 1	1.0714	1.0714	10194.00	11351.61	10000.000	-----
19	11.861 4-CHLOROCINE	1.0701	1.0701	10194.00	11351.61	10000.000	-----
20	12.046	4.7391	4.7391	10194.00	11351.61	-----	4.000e5
21	12.175	1.1393	1.1393	10194.00	11351.61	-----	4.000e5
22	12.311	1.1313	1.1313	10194.00	11351.61	-----	4.000e5
23	12.405 4-BENZYLIC	1.1311	1.1311	10194.00	11351.61	10000.000	-----
24	12.430 EDO	1.1303	1.1303	10194.00	11351.61	10000.000	-----
25	12.511	11.1533	11.1533	11.1533	11.1533	-----	4.000e5
26	13.011 ENOD 1	-0.1044	-0.1044	10194.00	11351.61	10000.000	-----
27	13.215 EDO 1	-0.1037	-0.1037	10194.00	11351.61	10000.000	-----
28	13.457	31.1762	31.1762	10194.00	11351.61	-----	4.000e5
29	13.805 DDO	5.1954	5.1954	10194.00	11351.61	10000.000	-----
30	13.971 ENODIN ALDE	1.7352	1.7352	10194.00	11351.61	10000.000	-----
31	14.151	37.3677	37.3677	10194.00	11351.61	-----	4.000e5
32	14.189 ENODOLATE	9.4441	9.4441	3037410.25	507456.91	10000.000	4.000e5
33	14.636 EOT	3.9817	3.9817	10194.00	11351.61	10000.000	4.000e5
34	14.597	31.2673	31.2673	10194.00	11351.61	-----	4.000e5
35	15.122	2.4403	2.4403	1196150.75	208953.83	-----	4.000e5
36	15.181	1.2163	1.2163	10194.00	11351.61	-----	4.000e5
37	15.522	0.5863	0.5863	587475.13	104815.27	-----	4.000e5
38	15.669 ENODIN KETONE	5.5063	5.5063	10194.00	11351.61	10000.000	4.000e5
39	16.111	0.8070	0.8070	10194.00	11351.61	-----	4.000e5

00022

1	1.011	-0.4531	101.0110	101.0110	-0.4531
2	1.012	-0.4531	101.0110	101.0110	-0.4531
3	10.013	10.0130	101.0110	101.0110	10.0130
4	10.014	-0.4531	101.0110	101.0110	-0.4531
5	10.015	10.0150	101.0110	101.0110	10.0150
6	10.016	-0.4531	101.0110	101.0110	-0.4531
7	10.017	10.0170	101.0110	101.0110	10.0170
8	10.018	-0.4531	101.0110	101.0110	-0.4531
9	10.019	10.0190	101.0110	101.0110	10.0190
10	10.020	-0.4531	101.0110	101.0110	-0.4531

RA.519 108.5119 1.010 1.010

Varian 3400 Capillary Column GC: 08-1701 30M x 0.53mm I.D. Detector: 3000C.
Injector: 150°C

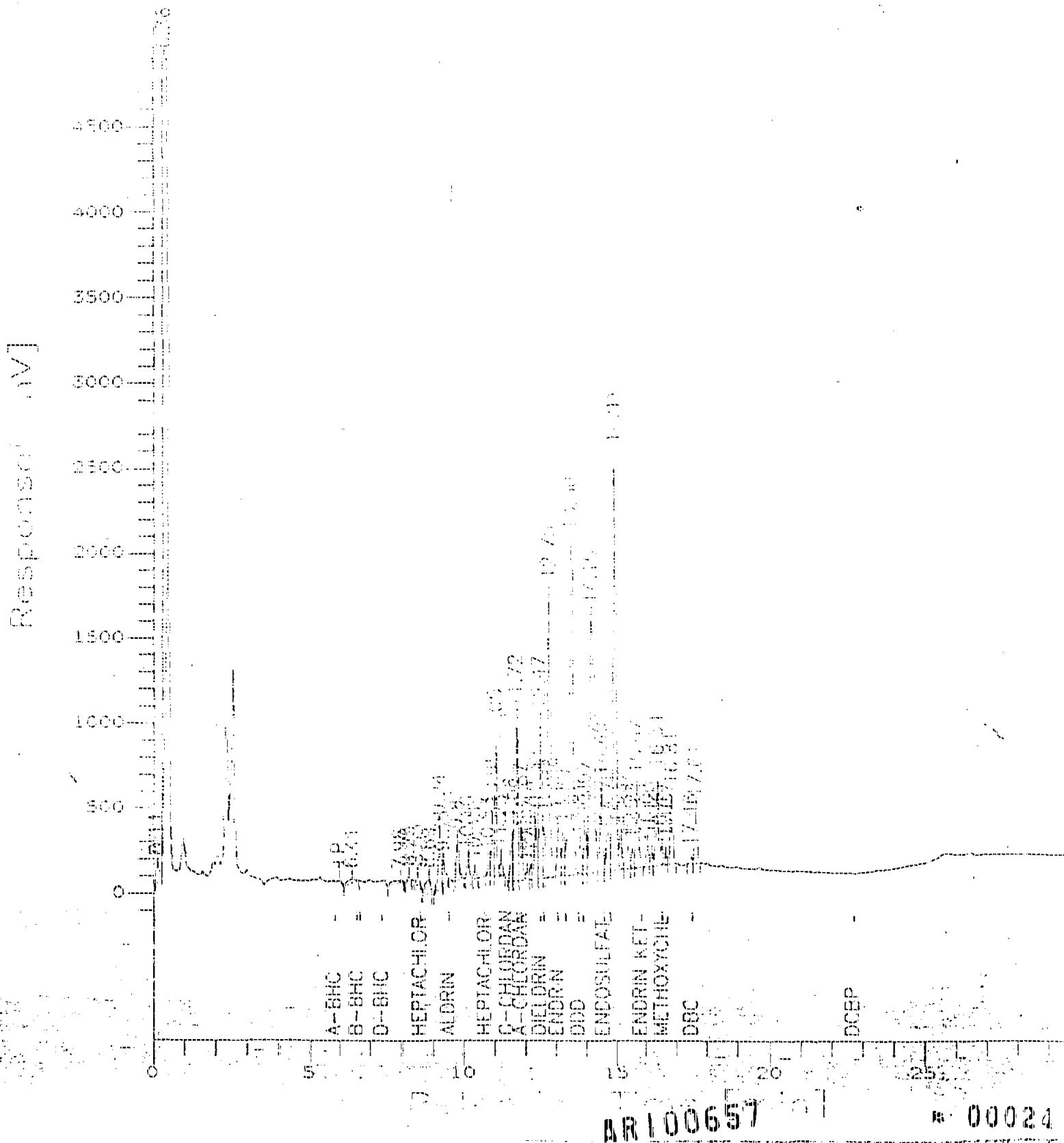
AR100656

00023

Retention Time (min) Chromatogram

Sample Name : 1004-00001
 Silencer : 1004-00001-00001.msp
 Start File : 0.00 min End File : 10.00 min
 Scale Factor : 1 Ret. Offset : 100.00

Sample ID : 1
 Date Sampled : 10/01/00
 Date Analyzed : 10/01/00
 Date Entered : 10/01/00
 Date Modified : 10/01/00



Software Version: 3.1 4000

Date: 10/24/91 17:09

Sample Name: 11260-EGCGA1

Data File: d:\2700\DATA\11264J01.D Date: 10/24/91 Job: 67

Sequence File: D:\2700\DATA\11264J01.seq File: 5 Channel: 1

Instrument: VARIAN 3100 Rack Vial: 170 Operator: JCB

Sample Amount: 1.0000 Dilution Factor: 1.0000

Princeton Testing Laboratory Analysis Report

PCPMTICRS AND PCP-1 BY CAPILLARY CHROMATOGRAPHY

Peak Ret Time t (min)	Concnen ppm	Rel. Peak Concentration (ppm)	Area (counts)	Height H (counts) (ppm)	Rel. Height H/ H ₁ (%)	Reference Retention	Rel. Calibra-
							tion Factor
1	0.114	0.004	000001.00	0.000000.00	-----		0.0000
2	0.114	0.100	000001.00	0.000000.00	-----		0.0000
3	0.143	-0.006	-0.0000	0.000000.00	-----		0.0000
4	0.213	0.500	000001.00	0.000000.00	-----		0.0000
5	11.035	HEPTACHLOR EPIC	0.8944	210456.00	418911.00	0.0	0.0000
6	11.515	3-CHLORDANE	0.7143	017401.00	192116.00	0.0	0.0000
7	11.702	ENDO 1	0.1151	179014.00	191207.00	-0.0000	0.0000
8	12.501		0.0000	0.000000.00	0.0000		0.0000
9	13.451	OIBLOBIN	0.0177	000001.00	145915.00	0.0	0.0000
10	13.483	DBE	0.0000	000001.00	112241.00	0.0	0.0000
11	13.724		0.0000	0.000000.00	0.0000		0.0000
12	13.974	ENDOTIN	0.5001	017955.00	1.0000	0.0	0.0000
13	13.978		0.0000	000001.00	0.000000.00	-----	0.0000
14	13.997	ENDO 1	0.0000	000001.00	1.0000	0.0	0.0000
15	13.999	DBE	0.0000	000001.00	0.000000.00	-----	0.0000
16	13.999	OIBLOBIN	0.4640	000001.00	0.000000.00	0.0000	0.0000
17	14.122		0.0000	0.000000.00	0.0000		0.0000
18	14.425		0.0000	0.000000.00	0.0000		0.0000
19	14.702	ENDOSULFATE	0.0000	000001.00	0.000000.00	0.0000	0.0000
20	14.887	DBE	0.0000	000001.00	0.000000.00	0.0000	0.0000
21	15.122		0.0000	0.000000.00	0.0000		0.0000
22	15.581		0.0000	000001.00	0.000000.00	0.0000	0.0000
23	15.492		0.0000	0.000000.00	0.000000.00		0.0000
24	15.495		0.0000	000001.00	0.000000.00		0.0000
25	15.739		0.0000	000001.00	0.000000.00		0.0000
26	15.802	ENDOTIN-3-ONE	0.0000	000001.00	1.0000	0.0	0.0000
27	15.835		0.0000	000001.00	0.000000.00		0.0000
28	15.874		0.0000	000001.00	0.000000.00		0.0000
29	15.944	METHONONICACID	0.0000	000001.00	0.000000.00	-0.0000	0.0000
30	15.952		0.0000	000001.00	0.000000.00		0.0000
31	16.731		006.7693	04211376.00	9.0000	0.0	0.0000
32	17.091		0.0000	000001.00	006879.00	0.0000	0.0000
33	17.594	DBE	58.4092	26043426.00	4.49e6	0.5359	0.0000
34	17.829		23.9770	10691670.00	1.76e6	0.0	0.0000
35	17.981		26.6559	11975415.00	1.90e6	0.0	0.0000
36	18.452		1.2173	542046.25	33904.00	0.00	0.0000
37	18.895		11.2979	550345.00	703015.15	0.0	0.0000
38	19.126		0.7406	520244.25	49817.00	0.00	0.0000
39	19.638		25.5701	01424524.00	1.36e6	0.00	0.0000

AR100658

00023

08/07/01 09:07:11 1168 1168

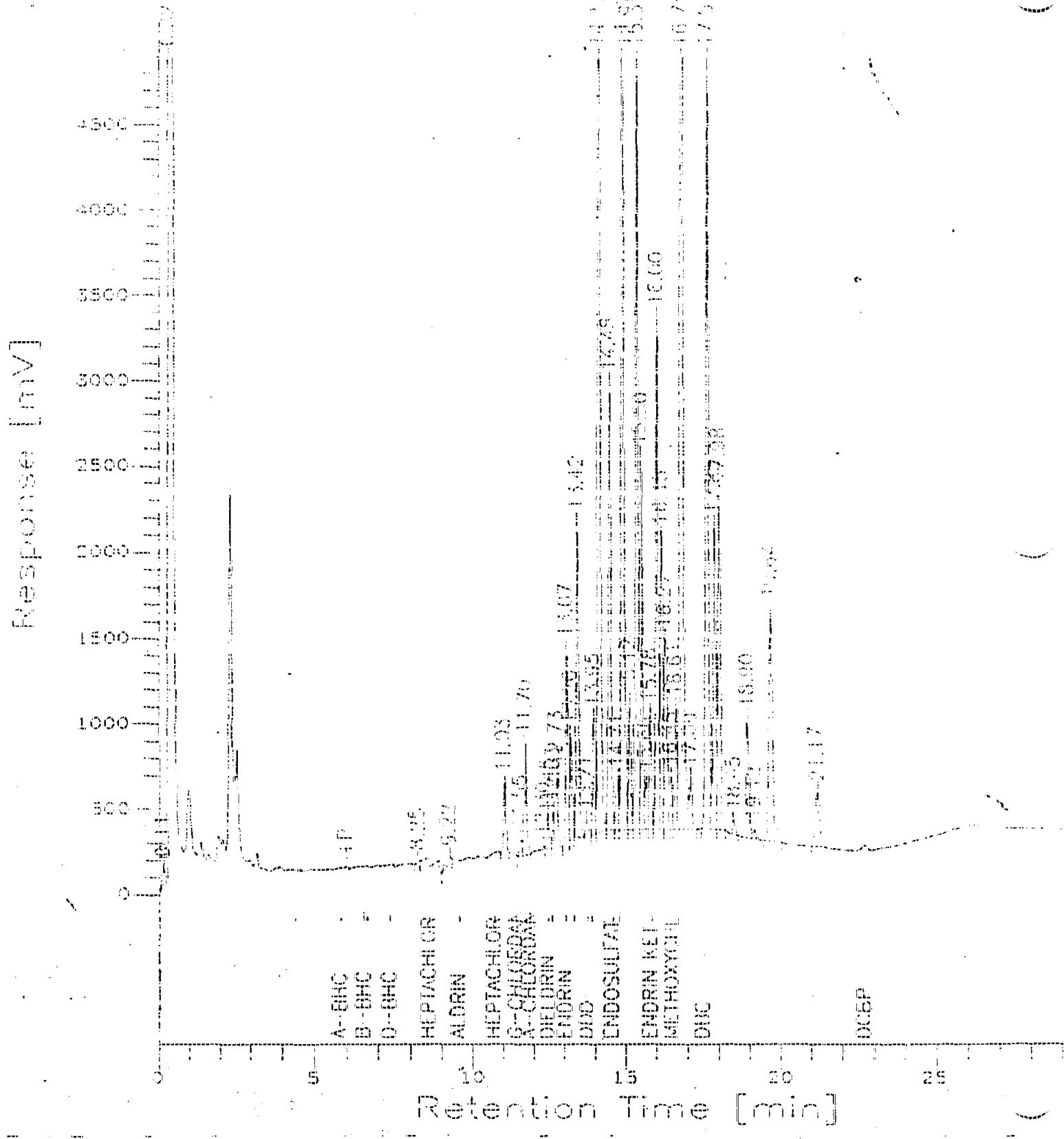
Varian 3400 Capillary Column GC: DB-1701 30m x 0.53mm I.D. 0.50um film Factor: 0.00 0

AR100659

* 00026

Baseline : 1000.00000
Release : 21.070000000000000
Start Time : 0.00 min End Time : 30.00 min
Scale Factor : 3 Plot Offset : -100.00

Baseline : 0
Date : 11/14/95 Page : 1
Log Start : 1000.000000 Log Point : 4000.000000
Plot Scale : 5000.00



**princeton testing
laboratory inc.**

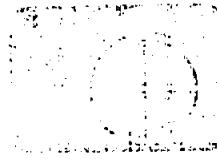
P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9050
FAX (609) 452-0347

Weston TAT
5 Underwood Court
Delran, New Jersey 08075
Attention: Marian Murphy

Report Date: 01/06/92
Job Number: 9108360-001GMWH
Date Received: 12/13/91

Analysis: Purgeable Volatile Organics By Method 8240 Page 1
Units: mg/kg

Parameters	Sample I.D.:	DF = 100,000	Dry Wgt. = 36.5%
	Sample D01		
Chloromethane		350J	
Bromomethane		<2,740	
Vinyl Chloride		<2,740	
Chloroethane		<2,740	
Methylene Chloride		1,620	
Acrolein		<13,700	
Acetone		<1,370	
Acrylonitrile		<13,700	
Carbon Disulfide		<1,370	
1,1-Dichloroethene		<1,370	
1,1-Dichloroethane		<1,370	
1,2-Dichloroethene (Total)		<1,370	
Trichlorofluoromethane		<1,370	
Chloroform		<1,370	
1,2-Dichloroethane		<1,370	
2-Butanone		384,000E 38%	
1,1,1-Trichloroethane		<1,370	
Carbon Tetrachloride		<1,370	
Vinyl Acetate		<1,370	
Bromodichloromethane		<1,370	
1,1,2,2-Tetrachloroethane		<1,370	
1,2-Dichloropropane		<1,370	
Trans-1,3-Dichloropropene		<1,370	
Trichlorosthene		<1,370	
Dibromochloromethane		<1,370	
1,1,2-Trichloroethane		<1,370	
Benzene		<1,370	
Cis-1,3-Dichloropropene		<1,370	
2-Chloroethylvinylether		<1,370	
Bromoform		<1,370	
2-Hexanone		<1,370	
4-Methyl-2-Pentanone		<1,370	
Tetrachloroethene		<1,370	
Toluene		160,600E 16%	
Chlorobenzene		<1,370	
Ethylbenzene		25,500	
Styrene		<1,370	
Total Xylenes		61,600	
1,3-Dichlorobenzene		<1,370	



princeton testing laboratory inc.

Weston TAT
5 Underwood Court
Delran, New Jersey 08075
Attention: Marian Murphy

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9
FAX (609) 452-0

Report Date: 01/06/92
Job Number: 9108360-001GMWH
Date Received: 12/13/91

Analysis: Purgeable Volatile Organics By Method 8240 Page 2

Units: $\mu\text{g}/\text{kg}$

DF = 100,000 Dry Wgt. = 36.5%

Parameters

Sample I.D.:

Sample D01

1,4-Dichlorobenzene

<1,370

1,2-Dichlorobenzene

<1,370

RECOVERY DATA

QC LIMITS

1,2-Dichloroethane-d4 (Surr.) 76-114 106

Toluene-d8 (Surr.) 88-110 95

4-Bromofluorobenzene (Surr.) 86-115 97

AR100662

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00029



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Princeton, NJ 08543-3108
(609) 452-9050
FAX (609) 452-0347

Weston TAT
5 Underwood Court
Delran, New Jersey 08075
Attention: Marian Murphy

Report Date: 01/06/92
Job Number: 9108360-001GMWH
Date Received: 12/13/91

Analysis: Volatile Organics By GCMS - Method 624 Page 1

Units: ug/ml

Parameters	Sample F.D.	DF = 10,000	Sample D05	DF = 10,000	Sample D06	Blank 12/20/91
Chloromethane		<100			12J	<10
Bromomethane		<100			<100	<10
Vinyl Chloride		<100			<100	<10
Chloroethane		<100			<100	<10
Methylene Chloride		88J			32J	2.6J
Acrolein		500			500	4.50
Acetone		580			330	4.4J
Acrylonitrile		500			500	50
Carbon Disulfide		500			500	50
1,1-Dichloroethene		500			500	50
1,1-Dichloroethane		500			500	50
1,2-Dichloroethene (Total)		500			500	50
Trichlorofluoromethane		500			500	50
Chloroform		500			500	50
1,2-Dichloroethane		500			500	50
2-Butanone		500			500	50
1,1,1-Trichloroethane		500			500	50
Carbon Tetrachloride		500			500	50
Vinyl Acetate		500			500	50
Bromodichloromethane		500			500	50
1,1,2,2-Tetrachloroethane		500			500	50
1,2-Dichloropropane		500			500	50
Trans-1,3-Dichloropropene		500			500	50
Trichloroethene		500			170	50
Dibromochloromethane		500			500	50
1,1,2-Trichloroethane		500			500	50
Benzene		500			210	50
Cis-1,3-Dichloropropene		500			500	50
2-Chloroethylvinylether		500			500	50
Bromoform		500			500	50
2-Hexanone		500			500	50
4-Methyl-2-Pentanone		500			500	50
Tetrachloroethene		500			23,700E	50
Toluene		560			560	50
Chlorobenzene		500			500	50
Ethylbenzene		7050			500	50
Styrene		450			500	50
Total Xylenes		13,300E			460	50
1,3-Dichlorobenzene		<100			<100	50

AR100663



princeton testing laboratory inc.

P.O. Box 3108
3490 U.S. Route 1
Princeton, NJ 08543-3108
(609) 452-9
FAX (609) 452-6

Weston TAT
5 Underwood Court
Delran, New Jersey 08075
Attention: Marian Murphy

Report Date: 01/06/92
Job Number: 9108360-001GMWH
Date Received: 12/13/91

Analysis: Volatile Organics By GCMS -- Method 624 Page 2

Units: $\mu\text{g}/\text{ml}$

Parameters	Sample I.D.:	DF = 10,000	DF = 10,000	Blank 12/20/91
1,4-Dichlorobenzene		<100	<100	<10
1,2-Dichlorobenzene		<100	<100	<10

RECOVERY DATA (% RECOVERY) & QC LIMITS (% RECOVERY)

1,2-Dichloroethane-d4 (Surr.)	76-114	118**	114	87
Toluene-d8 (Surr.)	88-110	105	105	103
4-Bromofluorobenzene (Surr.)	86-115	98	NR	103

E = Estimated value above linear range.

NR = No recovery data due to matrix interference.

J = Estimated Value Detected Below MDL
** = Outside Of QC Limits

Bipin Patel, Ph.D.
Organic Laboratory Manager

For inquiries call us at (609) 452-9050 and ask for our Customer Service Department

Member: American Council of Independent Laboratories, Inc.

00035

LABORATORY CHRONICLE
ORGANIC ANALYSIS

Company: Electra-Tek Job #: 240326-0

Date Received & Refrigerated: 13/02/91

EXTRACTION INFORMATION

Base-Neutral Extractables

____ / /
____ / /
____ / /
____ / /
____ / /

Acid Extractables

____ / /
____ / /
____ / /

Pesticides/ PCBs

____ / /
____ / /

PCBs only

____ / /
____ / /

Herbicides

____ / /
____ / /

Pesticides (EPTOX)

____ / /

Other:

____ / /

ANALYSIS INFORMATION

Base-Neutral Extractables

____ / /
____ / /
____ / /
____ / /
____ / /

Acid Extractables

____ / /
____ / /
____ / /

Pesticides/ PCBs

____ / /
____ / /

PCBs only

____ / /
____ / /

Herbicides

____ / /
____ / /

Pesticides (EPTOX)

____ / /

Volatiles - 601/602

____ / /
____ / /

Volatiles - 624/8240

CC 13/02/91
CS 1/02/91 - m:msd
____ / /

Other:

____ / /

Dept. Manager Review and Approval: _____

QC Supervisor Review and Approval: Michael H-J/S1

BR100665

00032

PRINCIPAL OCCURRENCE

Sample ID: 100666-001
 Sample Type: 100666-001
 Instrument File:
 Date to 8307 sample
 Lab Number:

Date: 07/31/89 at 23:56
 Calcd: CALTAG 0 3
 Analysis: GC

Page No.: 95
 File No.: 144364
 Job No.: 8363

Laboratory:

Contract:

m/z	Intensity	X RA	Ion Abundance Criteria			Actual	Status
			Min %	Max %	Mass		
80	50.02	17.8	15.0	40.0	95	17.6	PASS
79	12160.	43.1	33.0	60.0	95	43.1	PASS
78	13224.	100.0	100.0	----	----	100.0	PASS
76	2140.	7.6	5.0	9.0	95	7.5	PASS
173	15.	0.1	----	2.0	174	0.1	PASS
174	21536.	76.3	50.0	----	95	76.3	PASS
175	1526.	5.4	5.0	9.0	174	7.1	PASS
176	21120.	74.8	95.0	101.0	174	95.1	PASS
177	1374.	4.9	5.0	9.0	176	6.5	PASS

AR100666

v 0003

DATE RECEIVED IN FILE
 1967 NOV 10 1967 DATE TAKEN 1967 NOV 10
 PAGE SPECTROGRAPHIC TYPE CHECKED
 PLEASE PLACE BACK IN COLUMN-STRIP DRAWING BOXES
 AND DO NOT DESTROY

Line	O. CO	O.	Minima	Maxima	Min Intens.	% RA	O.	Intens.
Phase	X RA	Intens.		Mass		X RA		
36?	1. 11	313.		93		3. 25	917.	
37?	3. 26	1484.		94		9. 30	2624.	
38?	4. 54	1290.		95	100.00	29224.		
39?	3. 18	895.		96		7. 53	2140.	
40?	4. 13	1145.		97		0. 05	15.	
41?	2. 22	625.		98		0. 23	65.	
42?	1. 45	410.		103		0. 47	132.	
43?	5. 16	1455.		104		0. 07	20.	
44?	16. 10	4544.		105		0. 46	131.	
45?	10. 08	2844.		115		0. 24	68.	
46?	0. 06	17.		117		0. 90	254.	
47?	1. 47	414.		119		5. 65	1593.	
48?	0. 19	53.		120		0. 53	149.	
49?	3. 60	1017.		128		0. 11	31.	
50?	17. 83	5032.		130		0. 60	163.	
51?	5. 92	1672.		134		2. 50	705.	
52?	0. 09	24.		135		0. 15	43.	
53?	0. 07	21.		141		0. 22	63.	
55?	1. 58	447.		143		0. 16	44.	
56?	1. 99	561.		155		0. 05	15.	
57?	4. 39	1238.		173		0. 05	15.	
58?	3. 02	853.		174		76. 30	21505.	
59?	1. 59	449.		175		5. 41	1525.	
60?	0. 93	263.		176		74. 83	21120.	
61?	3. 76	1062.		177		4. 87	1374.	
62?	3. 42	966.		207		0. 95	244.	
63?	2. 70	763.						
65?	0. 07	21.						
67?	0. 44	125.						
68?	8. 87	2504.						
69	8. 80	2484.						
70	1. 24	350.						
71	0. 64	182.						
72	1. 19	336.						
73	6. 74	1702.						
74	12. 64	3568.						
75	43. 08	12160.						
76	3. 79	1065.						
77	1. 36	364.						
78	0. 54	151.						
79	1. 63	459.						
80	0. 19	55.						
81	1. 64	520.						
82	0. 61	172.						
86	0. 78	219.						
87	5. 18	1462.						
88	5. 39	1604.						
89	1. 74	492.						
91	1. 47	416.						
92	2. 07	584.						

AR100667

0003

MASS SPECTRUM
12/20/81 9:35:00 + 5:24
SAMPLE: 5046 SF MASS SPECTROMETER TUNE CHECK
COLUMNS: CLASSIC PACKED COLUMN/2METHYL 624/1.0 10/18/91
TEMP: 239 DEG. C
#255 10 #257 SUMMED

95

DATA: BFB1220 #255
CALL: GA-TEC #3
SIC: 141204.

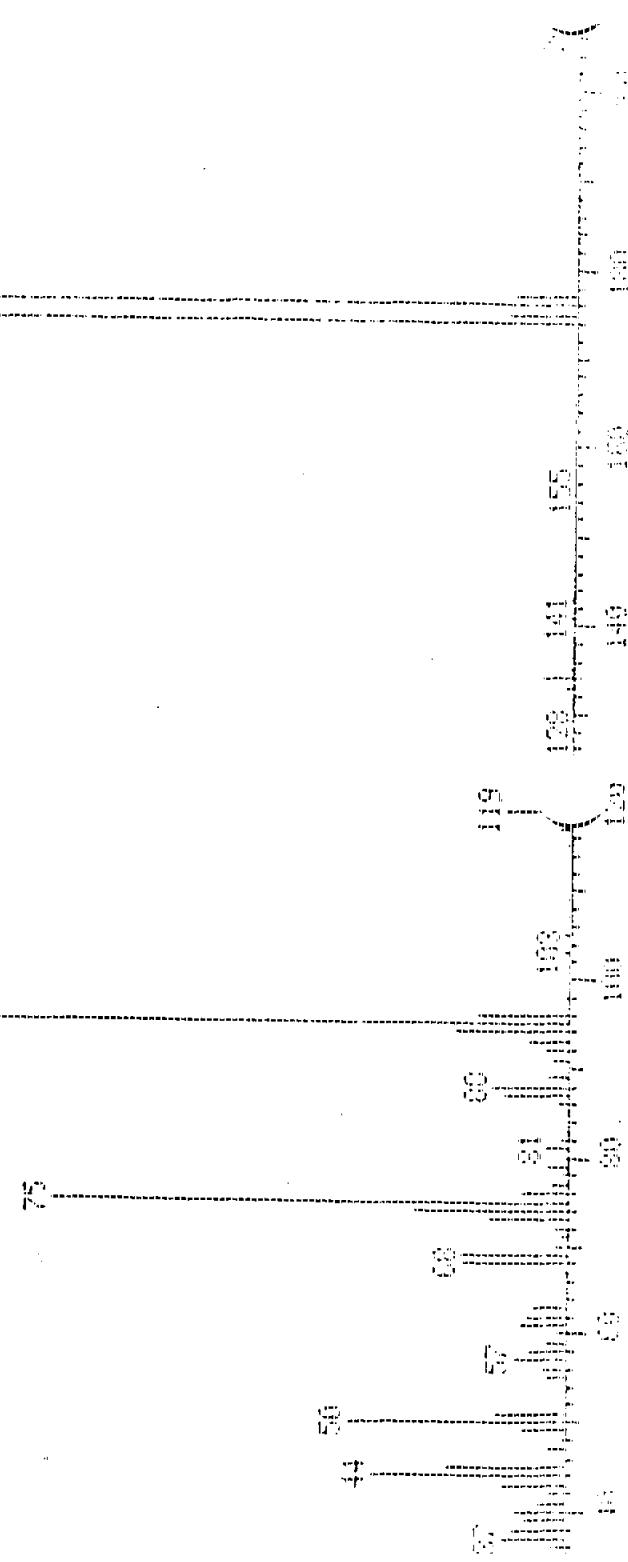
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H2: 141204.

100.0

50.0

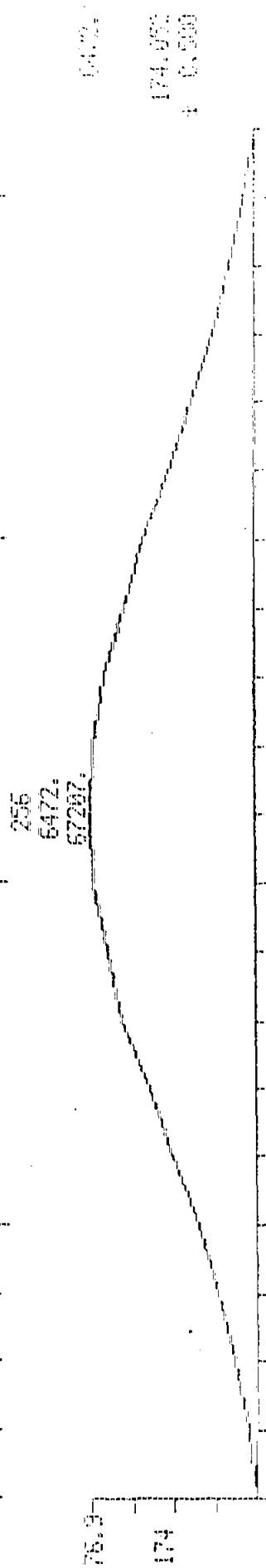
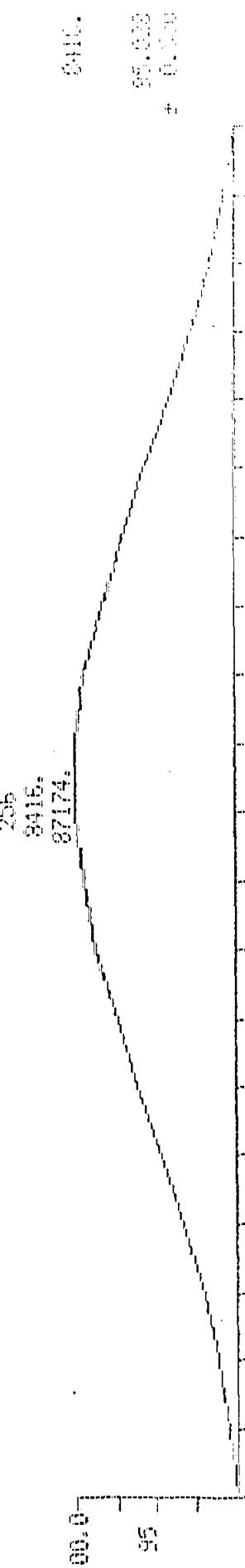
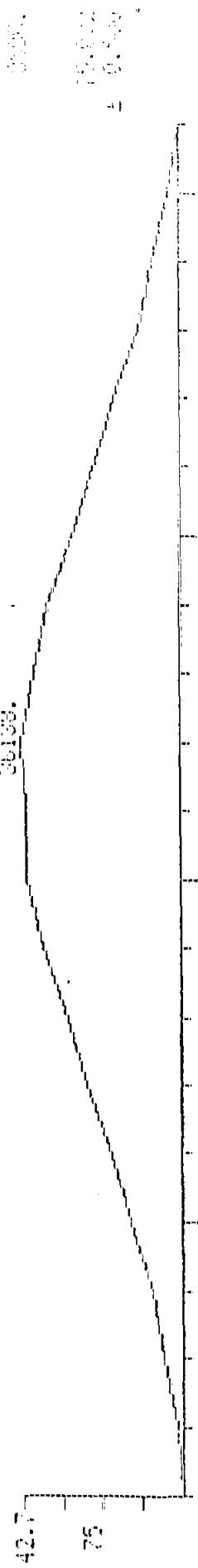
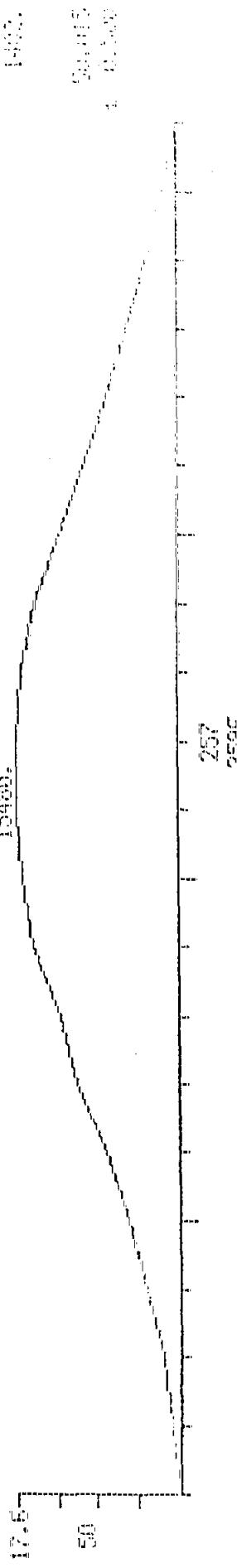
AB 100668

00035



MASS CHROMATOGRAMS
12/26/91 9:35:09

DATA: BFB1220 #255
CALI: CALTAB #3
SAMPLE: 50% BFB 40% SPECIOMETER TIME CHECK
CONC.: CLASS PACKED COLUMN/METHOD .624/1-C 16/18/91
PUMP: G 1, 400 LABEL: N 3, 4.0 QUANT: A 3, 1.02 J 0 BASE: U 20, 3



BR/00669

00036

CHLOROFLUORODENZENE

100-6773

Testing Report
 Date: 02/22/00 + 5:20
 Instrument: FINN
 File No: 8477-
 Date Number:

Date: 02/21/00 # 265
 Calib: CALTAG # 3
 Analyst: CC

Base m/z: 95
 RIC: 211200.
 Acct. No.: 8477

Laboratory:

Contract:

m/z	Intensity	% RA	Ion Abundance Criteria		Mass	Actual	Status
			Min %	Max %			
50	6175.	15.7	15.0	40.0	95	15.7	PASS
75	14735.	37.6	30.0	60.0	95	37.6	PASS
95	39232.	100.0	100.0	----	95	100.0	PASS
95	2868.	7.3	5.0	9.0	95	7.3	PASS
173	174.	0.4	----	2.0	174	0.4	PASS
174	27072.	69.0	50.0	----	95	69.0	PASS
175	1934.	4.9	5.0	9.0	174	7.1	PASS
176	25952.	66.2	95.0	101.0	174	95.9	PASS
177	1754.	4.5	5.0	9.0	176	6.8	PASS

AR100670

0003

Mass List

01/02/72 8:32:00 + 6:23

Sample: 5049-BFB

Conc.: CLASS PACKED COLUMN METHOD A24

#254 50-#253 summed

Notes: DB12E + 255

Col1: CALTAG # 2

Mass SPECTROMETER TUNE CHECK

Date m/z: 93

Run: 211200.

S4 250	0.00	0	Minima	Min Inten	ATR
			Maxima	at 0	
34?	0. 00	0	Mass	% RA	Inten
35?	0. 92	361.	89	2. 95	3120.
37?	4. 89	1920.	89	2. 48	971.
38?	4. 45	1744.	91	1. 05	413.
39?	4. 14	1626.	92	1. 78	700.
40?	4. 06	1592.	93	2. 99	1172.
41?	6. 28	2464.	94	8. 42	3304.
42?	2. 89	1134.	95	100. 00	39232.
43?	9. 24	3624.	96	7. 31	2868.
44?	19. 17	7520.	97	1. 03	404.
45?	15. 52	6088.	98	0. 49	194.
46?	0. 44	174.	101	0. 34	134.
47?	1. 95	764.	102	0. 42	163.
48?	0. 43	169.	103	0. 66	259.
49?	3. 15	1236.	104	0. 09	34.
50?	15. 74	6176.	105	0. 79	311.
51?	5. 12	2008.	106	0. 20	77.
52?	0. 29	1113.	107	0. 15	60.
53?	0. 76	297.	109	0. 11	45.
54?	0. 51	202.	110	0. 32	124.
55?	5. 44	2136.	111	0. 41	160.
56?	3. 84	1508.	112	1. 51	593.
57?	8. 64	3888.	113	0. 50	196.
58?	4. 03	1580.	115	0. 65	255.
59?	2. 38	903.	116	0. 24	93.
60?	1. 36	532.	117	0. 95	374.
61?	3. 34	1312.	118	0. 30	117.
62?	2. 99	1174.	119	1. 05	413.
63?	2. 29	897.	125	0. 19	74.
65?	0. 27	106.	128	0. 72	282.
67?	1. 29	507.	129	0. 80	312.
68?	7. 71	3024.	130	0. 39	152.
69	9. 42	3696.	131	0. 75	293.
70	3. 59	1410.	133	0. 99	390.
71	2. 92	1146.	141	0. 51	202.
72	2. 09	820.	143	0. 92	361.
73	6. 27	2460.	145	0. 98	393.
74	10. 87	4264.	147	0. 43	167.
75	37. 56	14736.	152	0. 12	46.
76	3. 63	1426.	157	0. 40	155.
77	1. 33	523.	159	0. 52	204.
78	0. 59	233.	171	1. 88	739.
79	1. 76	692.	172	0. 16	63.
80	0. 56	218.	173	0. 44	174.
81	2. 12	830.	174	69. 00	27072.
82	1. 01	398.	175	4. 93	1934.
83	1. 87	732.	176	66. 15	25952.
84	1. 67	655.	177	4. 47	1754.
85	0. 75	295.	205	1. 42	556.
86	1. 26	493.	207	0. 36	143.
87	6. 87	2696.	220	0. 22	85.

AR100671

00038

MS SPECTRUM
91/02/22 8:32:00 + 5:23
SAMPLE: 5015 BPC MASS SPECTROMETER TIME CHECK
00005.1 GLASS PACKED COLUMN/HEAT 624

TEMP: 226 DEG. C
#254 10 #255 SWINED

95

DATE: 07/01/2 #255
CALL: CALTAG #2

BASE M/Z: 95
PIC: 211265.

100.0

200.0

1

100

70

80

90

100

110

120

130

140

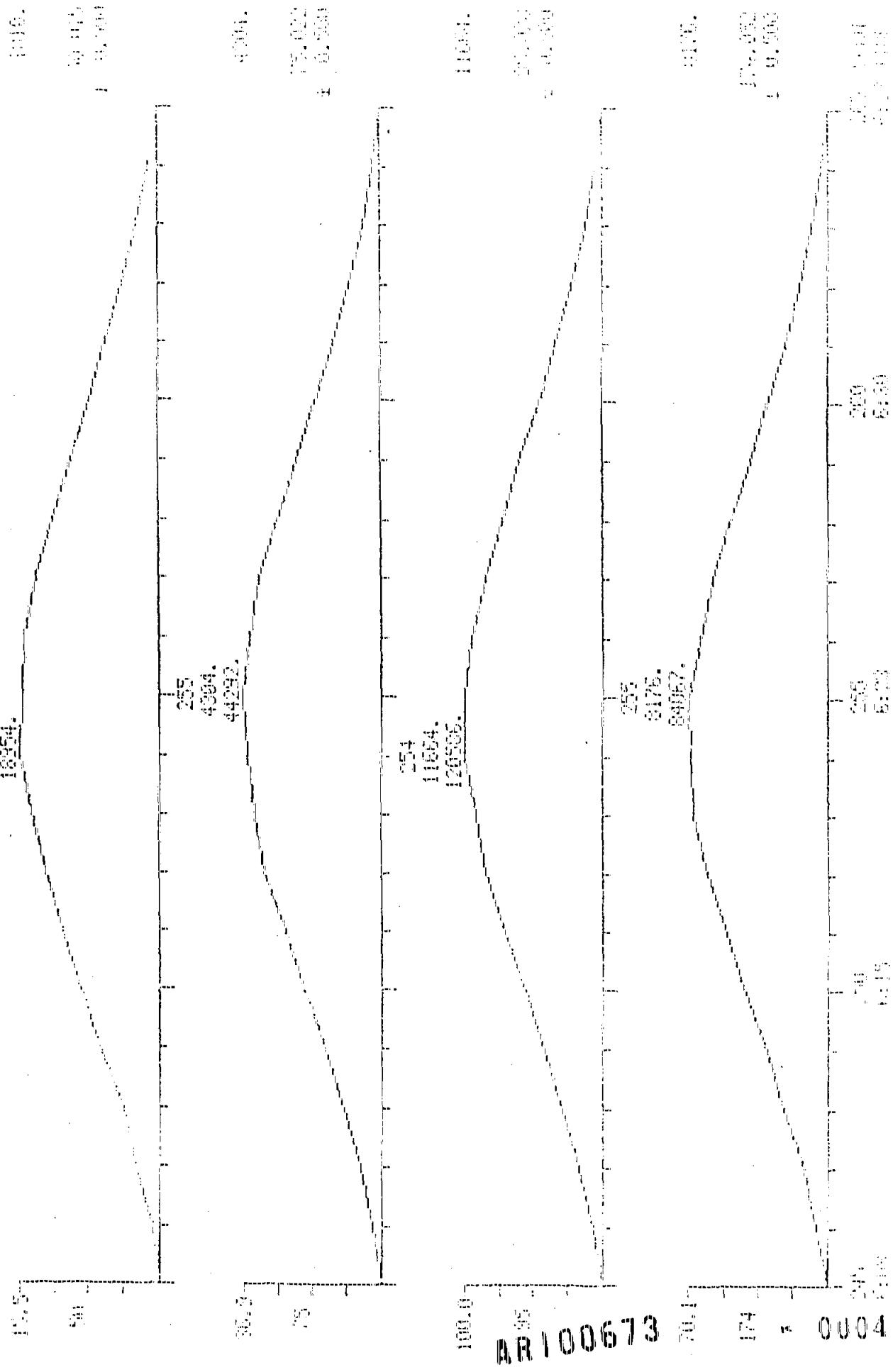
AR 100672

• 00030



MESS CHROMATOGRAMS
01/02/92 8:32:00
SAMPLE: 5496 LFB MASS SPECTROMETER TUNE CHECK
COMBINE: 01/02/92 00:00:00
PATTERN: 0 1, 400, LABEL: H 3, 4.0 QMIN: 0.3, 1.6 Q 0 BASE: 0 20, 3

DATA: EFB12 #254
CAL1: CALTEC 43
SCANS: 245 TO 255



20100613 0004

Initial Calibration Data

Instrument Identifier: FINI

Calibration Date 12-18-91

MAX AVG. XRS FOR CCC = .30

MIN AVG. RF FOR SPCC 0.300

Compound	RF 10 NG	RF 20 NG	RF 50 NG	RF 100NG	RF 200NG	Mean RF	XRS
C010 CHLOROMETHANE **	0. 947	0. 857	0. 949	0. 812	0. 903	0. 874	6. 62
C015 BROMOMETHANE	0. 764	0. 651	0. 484	0. 238	0. 181	0. 464	54. 61
C020 VINYL CHLORIDE *	0. 607	0. 573	0. 620	0. 545	0. 501	0. 529	24. 74
C025 CHLOROETHANE	0. 296	0. 295	0. 276	0. 221	0. 262	0. 270	11. 91
C030 METHYLENE CHLORIDE	1. 036	1. 003	0. 921	0. 736	0. 710	0. 891	17. 11
C031 ACROLIN	0. 227	0. 271	0. 274	0. 268	0. 324	0. 273	12. 55
C035 ACETONE	0. 177	0. 185	0. 133	0. 127	0. 138	0. 152	17. 61
C052 ACRYLONITRILE	0. 550	0. 660	0. 635	0. 664	0. 823	0. 666	14. 81
C040 CARBON DISULFIDE *	1. 612	1. 631	1. 893	1. 717	1. 756	1. 722	6. 51
C045 1,1-DICHLOROETHENE **	0. 614	0. 612	0. 759	0. 701	0. 775	0. 692	11. 18
C050 1,1-DICHLOROETHANE **	1. 331	1. 399	1. 593	1. 542	1. 720	1. 517	10. 27
C055 TRANS-1,2-DICHLOROETHEN	0. 974	1. 025	1. 183	1. 170	1. 321	1. 134	12. 17
C060 TRICHLOROFLUOROMETHA	1. 328	1. 304	1. 280	1. 286	0. 869	1. 213	151. 9
C065 CHLOROFORM *	1. 668	1. 607	1. 782	1. 709	1. 936	1. 740	7. 21
C066 1,2-DICHLOROETHANE	0. 876	0. 839	0. 964	0. 954	1. 058	0. 938	9. 08
C110 2-BUTANONE	0. 035	0. 057	0. 053	0. 053	0. 058	0. 051	17. 91
C115 1,1,1-TRICHLOROETHANE	0. 541	0. 553	0. 591	0. 556	0. 575	0. 563	4. 41
C120 CARBON TETRACHLORIDE	0. 436	0. 440	0. 528	0. 486	0. 517	0. 481	7. 71
C125 VINYL ACETATE	0. 179	0. 192	0. 258	0. 194	0. 264	0. 217	10. 57
C130 BROMO DICHLOROMETHANE	0. 543	0. 561	0. 584	0. 587	0. 606	0. 576	4. 81
C140 1,2-DICHLOROPROPANE	0. 321	0. 345	0. 353	0. 371	0. 359	0. 350	5. 21
C145 TRANS-1,3-DICHLOROPROPE	0. 695	0. 828	0. 798	0. 844	0. 860	0. 805	8. 1
C150 TRICHLOROETHENE	0. 382	0. 417	0. 436	0. 422	0. 425	0. 416	4. 9
C155 DIBROMOCHLOROMETHANE	0. 478	0. 508	0. 545	0. 552	0. 565	0. 529	6. 7
C160 1,1,2-TRICHLOROETHANE	0. 291	0. 303	0. 302	0. 318	0. 320	0. 305	5. 1
C165 BENZENE	0. 862	0. 873	0. 897	0. 905	0. 912	0. 890	2. 4
C143 CIS-1,3-DICHLOROPROPENE	0. 695	0. 828	0. 798	0. 844	0. 860	0. 805	8. 1
C175 2-CHLOROETHYL VINYL ETH	0. 146	0. 099	0. 164	0. 183	0. 193	0. 157	23. 6
C180 BROMOFORM **	0. 343	0. 390	0. 379	0. 406	0. 415	0. 388	6. 7
C220 TETRACHLOROETHENE	0. 582	0. 621	0. 623	0. 595	0. 628	0. 610	3. 2
C210 2-HEXANONE	0. 120	0. 167	0. 124	0. 114	0. 130	0. 131	15. 9
C205 4-METHYL 2-PENTANONE	0. 192	0. 288	0. 231	0. 254	0. 261	0. 243	14. 4
C225 1,1,2,2-TETRACHLOROETHA	0. 370	0. 493	0. 422	0. 474	0. 467	0. 445	11. 1
C230 TOLUENE *	1. 126	1. 174	1. 174	1. 150	1. 160	1. 157	1. 7
C225 CHLOROBENZENE **	0. 947	0. 998	0. 958	0. 944	0. 959	0. 961	2. 2
C240 ETHYL BENZENE *	0. 441	0. 450	0. 444	0. 443	0. 454	0. 445	1. 1
C245 STYRENE	0. 844	0. 878	0. 864	0. 867	0. 891	0. 869	2. 0
C250 M-XYLENE	0. 506	0. 543	0. 500	0. 499	0. 540	0. 517	4. 2
C253 1,3-DICHLOROBENZENE	0. 506	0. 605	0. 520	0. 557	0. 555	0. 548	7. 0
C254 1,4-DICHLOROBENZENE	0. 497	0. 591	0. 455	0. 522	0. 515	0. 516	9. 5
C255 1,2-DICHLOROBENZENE	0. 615	0. 720	0. 681	0. 652	0. 695	0. 670	6. 0
C250 o,p-XYLENE	0. 974	1. 063	0. 990	0. 974	1. 024	1. 006	3. 7

AR100674

00041

W:\ECL\DATAFILE\SVO1.EOR

Calibration Check

Instrument Identifier: FINN

Calibration Date: 12-19-91

Standard File: SVO1.EOR

Date: 12/20/91 Time: 12:00:00

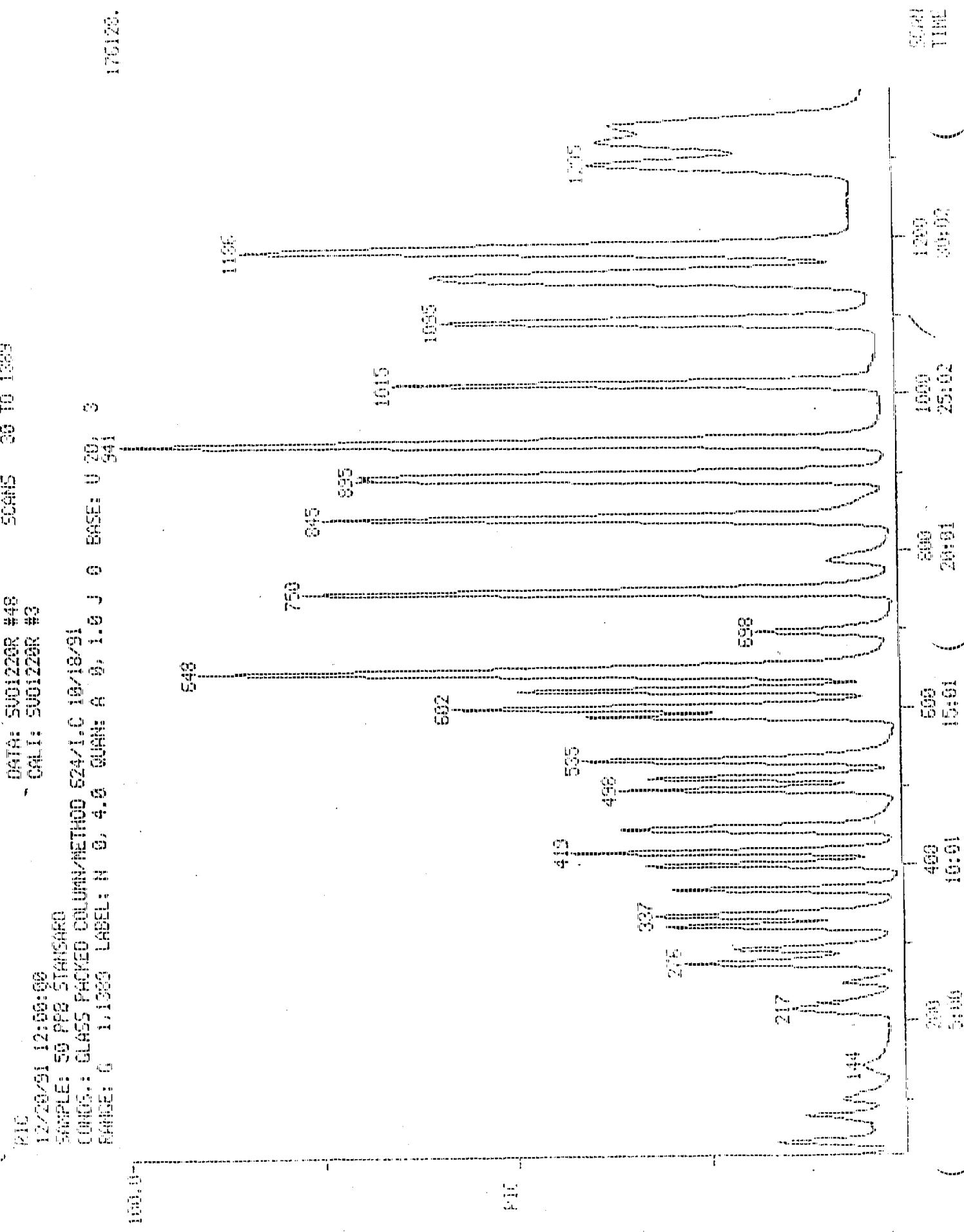
MAX RPD FOR CCC (*) = 25 %

MIN RPD FOR SPCC (**) = 0.300

Compound	Mean	RF(1)	RF(0)	Z.D.
C010 CHLOROMETHANE	**	0. 874	0. 946	5. 969
C015 BROMOMETHANE		0. 464	0. 443	3. 433
C020 VINYL CHLORIDE	*	0. 529	0. 562	5. 110
C025 CHLOROETHANE		0. 270	0. 325	20. 490
C030 METHYLENE CHLORIDE		0. 881	0. 670	24. 005
C031 ACROLIN		0. 273	0. 269	1. 174
C035 ACETONE		0. 152	0. 136	10. 180
C052 ACRYLONITRILE		0. 666	0. 706	6. 005
C040 CARBON DISULFIDE		1. 722	1. 620	5. 905
C045 1,1-DICHLOROETHENE	*	0. 692	0. 652	5. 824
C050 1,1-DICHLOROETHANE	**	1. 517	1. 464	3. 458
C055 TRANS-1,2-DICHLOROETHEN		1. 134	1. 053	7. 145
C060 TRICHLOROFLUOROMETHA		1. 213	0. 914	24. 686
C065 CHLOROFORM	*	1. 740	1. 725	0. 902
C085 1,2-DICHLOROETHANE		0. 938	0. 900	4. 085
C110 2-BUTANONE		0. 051	0. 049	3. 195
C115 1,1,1-TRICHLOROETHANE		0. 563	0. 504	10. 536
C120 CARBON TETRACHLORIDE		0. 481	0. 417	13. 289
C125 VINYL ACETATE		0. 217	0. 293	34. 755
C130 BROMO DICHLOROMETHANE		0. 576	0. 573	0. 530
C140 1,2-DICHLOROPROPANE		0. 350	0. 335	4. 334
C145 TRANS-1,3-DICHLOROPROPE		0. 805	0. 831	3. 198
C150 TRICHLOROETHENE		0. 416	0. 374	10. 205
C155 DIBROMOCHLOROMETHANE		0. 529	0. 533	0. 621
C160 1,1,2-TRICHLOROETHANE		0. 305	0. 320	4. 969
C165 BENZENE		0. 890	0. 908	2. 005
C143 CIS-1,3-DICHLOROPROPENE		0. 805	0. 831	3. 193
C175 2-CHLOROETHYL VINYL ET	H	0. 157	0. 188	19. 540
C180 BROMOFORM	**	0. 388	0. 402	3. 856
C220 TETRACHLOROETHENE		0. 610	0. 554	7. 426
C210 2-HEXANONE		0. 131	0. 111	15. 711
C205 4-METHYL 2-PENTANONE		0. 243	0. 224	7. 751
C225 1,1,2,2-TETRACHLOROETHA		0. 445	0. 464	4. 177
C230 TOLUENE	*	1. 157	1. 156	0. 035
C235 CHLOROBENZENE	**	0. 961	0. 961	0. 004
C240 ETHYL BENZENE	*	0. 446	0. 430	3. 688
C245 STYRENE		0. 869	0. 852	1. 929
C250 M-XYLENE		0. 517	0. 501	3. 104
C253 1,3-DICHLOROBENZENE		0. 548	0. 508	7. 370
C254 1,4-DICHLOROBENZENE		0. 516	0. 478	7. 359
C255 1,2-DICHLOROBENZENE		0. 673	0. 630	6. 447
C250 O&P XYLENE		1. 005	0. 974	3. 032

AR100675

00042



AR100676

• 00043

✓
excellen

Quantitation Report File: SV0128CR

Sample: SV0128CR

12-NOV-91 12:00:00

Serial: 50 PTFE STANDARD

Column: GLASS PACKED COLUMN/METHOD 6244 L.C. 10/18/91

Formula: STD

Instrument: FINI

Weight: 0.000

Submitted by PTL

Analyst: CC

Acct. No.: PTL

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fact. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **INT. STD. **
2	C110 1, 4-DIFLUOROBENZENE **INT. STD. **
3	C120 CHLOROBENZENE-D5 **INT. STD. **
4	C215 1, 2-DICHLOROETHANE-D4 **S. STD. **
5	C205 TOLUENE-D9 **S. STD. **
6	C210 4-BROMOFLUOROBENZENE **S. STD. **
7	C010 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1, 1-DICHLOROETHENE *
17	C050 1, 1-DICHLOROETHANE **
18	C055 TRANS-1, 2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1, 2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1, 1, 1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1, 2-DICHLOROPROPANE *
28	C145 TRANS-1, 3-DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1, 1, 2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1, 3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1, 1, 2, 2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M-XYLENE
45	C253 1, 3-DICHLOROBENZENE
46	C254 1, 4-DICHLOROBENZENE
47	C255 1, 2-DICHLOROBENZENE

AR100677

00044

DATAFILE: SV012R

Calibration Check

Instrument Identifier: FINN

Calibration Date: 12-13-91

Standard File: SV012R

Date: 01/02/92 Time: 11.08:00

MAX RPD FOR CCC (*) = 25 %

MIN RF FOR SPCC (**) = 0.300

Compound	Mean RF (1)	RF (0)	% D
C010 CHLOROMETHANE **	0.894	1.045	16.980
C015 BROMOMETHANE	0.464	0.457	0.711
C020 VINYL CHLORIDE *	0.529	0.574	8.468 ✓
C025 CHLOROETHANE	0.270	0.327	21.180
C030 METHYLENE CHLORIDE	0.881	0.800	9.185
C261 ACROLIN	0.273	0.234	14.021
C265 ACETONE	0.152	0.161	5.706
C262 ACRYLONITRILE	0.655	0.671	0.745
C040 CARBON DISULFIDE	1.722	1.926	11.869
C045 1,1-DICHLOROETHENE *	0.692	0.777	12.299 ✓
C050 1,1-DICHLOROETHANE **	1.517	1.521 ✓	0.313
C055 TRANS-1,2-DICHLOROETHEN	1.134	1.145	0.927
C000 TRICHLOROFLUOROMETHA	1.213	0.961	20.742
C060 CHLOROFORM *	1.740	1.543	11.334 ✓
C065 1,2-DICHLOROETHANE	0.938	0.671	28.530
C110 2-BUTANONE	0.051	0.047	8.866
C115 1,1,1-TRICHLOROETHANE	0.563	0.411	27.128
C120 CARBON TETRACHLORIDE	0.481	0.348	27.727
C125 VINYL ACETATE	0.217	0.220	1.174
C130 BROMO DICHLOROMETHANE	0.576	0.476	17.335
C140 1,2-DICHLOROPROPANE	0.350	0.370	5.616 ✓
C145 TRANS-1,3-DICHLOROPROPENE	0.805	0.785	2.459
C150 TRICHLOROETHENE	0.416	0.380	8.824
C155 DIBROMOCHLOROMETHANE	0.529	0.452	14.621
C160 1,1,2-TRICHLOROETHANE	0.305	0.302	0.877
C165 BENZENE	0.890	1.005	12.975
C143 CIS-1,3-DICHLOROPROPENE	0.805	0.785	2.459
C175 2-CHLOROETHYL VINYL ETHER	0.157	0.090	42.440
C180 BROMOFORM **	0.388	0.304 ✓	21.639
C220 TETRACHLOROETHENE	0.610	0.598	1.881
C210 2-HEXANONE	0.131	0.099	24.713
C205 4-METHYL 2-PENTANONE	0.243	0.219	9.805
C225 1,1,2,2-TETRACHLOROETHA	0.445	0.468 ✓	5.246
C230 TOLUENE *	1.157	1.247	7.750 ✓
C235 CHLOROBENZENE **	0.961	1.000 ✓	4.026
C240 ETHYL BENZENE *	0.446	0.443	0.885 ✓
C245 STYRENE	0.859	0.834	8.991
C250 M-XYLENE	0.517	0.486	6.143
C253 1,3-DICHLOROBENZENE	0.546	0.472	13.980
C254 1,4-DICHLOROBENZENE	0.516	0.431	16.470
C255 1,2-DICHLOROBENZENE	0.673	0.529	21.833
C260 O,p-XYLENE	1.005	0.960	4.423

AR100678

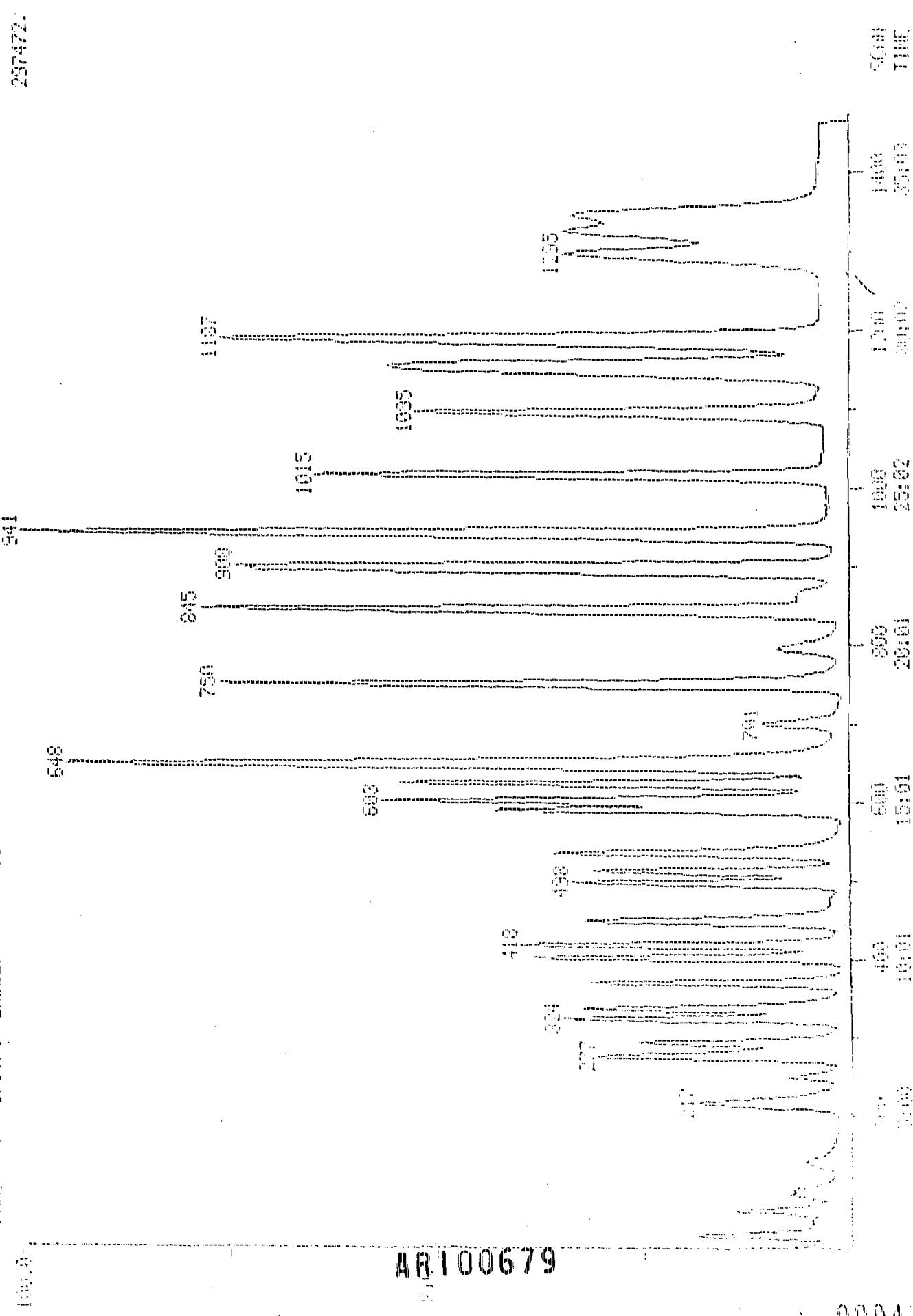
00046

PIC 01002732 11:08:49
SAMPLE 90 PPG ST6
POINT 2 01005 900750 624
PANEL 0 1,1474, 14681; 0, 4,9, 8,9, A, 0, 1,0, J, 0, 8,95, 0, 26, C
94, 1

SCA15 36 10 1474

DATA 200128 448
CALI 200128 43

207472.



AR100679

00047

Quantitation Report File: SVO1ER

Date: SVO1ER.TI

01/02/92 11:08:00

P.L.: 50 PPB STD

Col.: GLASS PACKED COLUMN/METHOD 624

Formula: SML

Instrument: FINN

Weight: 0.000

Submitted by: PTL

Analyst: CC

Acct. No.: PTL

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)
Resp. fact. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **INT. STD. **
2	C110 1, 4-DIFLUOROBENZENE **INT. STD. **
3	C120 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1, 2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	CO10 CHLOROMETHANE **
8	CO15 BROMOMETHANE
9	CO20 VINYL CHLORIDE *
10	CO25 CHLOROETHANE
11	CO30 METHYLENE CHLORIDE
12	C251 ACROLIN
13	CO35 ACETONE
14	C252 ACRYLONITRILE
15	CO40 CARBON DISULFIDE
16	CO45 1, 1-DICHLOROETHENE *
17	CO50 1, 1-DICHLOROETHANE **
18	CO55 TRANS-1, 2-DICHLOROETHENE
19	CO00 TRICHLOROFLUOROMETHANE
20	CO60 CHLOROFORM *
21	CO65 1, 2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1, 1, 1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1, 2-DICHLOROPROPANE *
28	C145 TRANS-1, 3-DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1, 1, 2-TRICHLOROETHANE
32	C165 BENZENE
33	C143 CIS-1, 3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1, 1, 2, 2-TETRACHLOROETHANE **
40	CE00 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M-XYLENE
45	C253 1, 3-DICHLOROBENZENE
46	C254 1, 4-DICHLOROBENZENE
47	C255 1, 2-DICHLOROBENZENE

BR100680

00045

No Name
42 C250 OSP XYLENE

No	m/z	Scan	Time	Ref	RRT	Meth	Area(Hight)	Amount	X Tot
1	49	337	6:26	1	1.000	A BB	225355.	50.000	1.85
2	114	749	18:45	2	1.000	A BB	599206	50.000	1.85
3	117	938	23:29	3	1.000	A BB	511161.	50.000	1.85
4	65	447	11:11	1	1.326	A BB	141310.	52.211	1.65
5	95	824	22:23	6	0.953	A BB	561311.	50.923	1.93
6	95	1035	27:24	3	1.167	A BB	372868.	50.423	1.87
7	50	47	1:11	1	0.139	A BB	235527.	51.217	1.84
8	94	79	1:59	1	0.234	A BB	105171.	53.599	1.98
9	62	101	2:32	1	0.300	M XX	129264.	57.901	2.14
10	64	142	3:33	1	0.421	A BB	73538.	50.731	1.88
11	49	217	5:26	1	0.644	A BB	180237.	48.754	1.80
12	56	251	6:17	1	0.745	A BB	52721.	40.208	1.49
13	43	250	6:15	1	0.742	A BB	36073.	42.738	1.58
14	53	275	6:53	1	0.816	A BB	151170.	47.740	1.77
15	75	277	6:56	1	0.822	A BB	434019.	50.138	1.85
16	96	324	8:07	1	0.961	A BB	175096.	60.768	2.25
17	63	370	9:16	1	1.078	A BB	342756.	65.085	2.41
18	61	402	10:04	1	1.193	A BB	257916.	51.925	1.92
19	101	294	7:22	1	0.872	A BB	216566.	109.529	4.05
20	83	418	10:28	1	1.240	A BB	347598.	59.696	2.21
21	62	452	11:19	1	1.341	A BB	151009.	56.395	2.05
22	43	477	11:56	2	0.637	A BB	27610.	63.000	2.33
23	97	498	12:28	2	0.665	A BB	245678.	74.901	2.77
24	117	512	12:49	2	0.684	A VB	208205.	76.520	2.83
25	43	537	13:26	2	0.717	A BB	131453.	75.525	2.79
26	83	536	13:25	2	0.716	A BB	285142.	55.116	2.04
27	63	591	14:48	2	0.789	A BB	221113.	56.151	2.08
28	75	602	15:04	2	0.804	A BB	470242.	56.905	2.10
29	95	624	15:37	2	0.833	A BB	227220.	54.594	2.02
30	129	644	16:07	2	0.860	A BB	270578.	52.871	1.96
31	97	652	16:19	2	0.870	A BB	180713.	53.567	1.98
32	78	649	16:13	2	0.865	A BB	602085.	53.014	1.96
33	75	602	15:04	2	0.804	A BB	470242.	56.905	2.10
34	63	700	17:31	2	0.935	A BB	53892.	52.719	1.95
35	173	751	18:48	2	1.003	A BB	181654.	52.023	1.92
36	166	845	21:09	3	0.901	A BB	305593.	51.808	1.92
37	43	862	21:35	3	0.919	A BB	50269.	52.573	1.94
38	43	796	19:56	3	0.849	A BB	111875.	58.251	2.15
39	83	840	21:02	3	0.896	A BB	239198.	55.363	2.05
40	91	901	22:33	3	0.961	A BB	636909.	52.285	1.93
41	112	943	23:36	3	1.005	A BB	510973.	51.545	1.91
42	106	1015	25:24	3	1.082	A BB	225937.	51.113	1.69
43	104	1147	28:43	3	1.223	A BB	426180.	52.087	1.93
44	106	1155	28:55	3	1.231	A BB	247987.	50.600	1.87
45	146	1294	32:23	3	1.380	A BB	240953.	54.336	2.01
46	146	1324	33:09	3	1.412	M XX	220048.	54.061	2.00
47	146	1342	33:35	3	1.431	M XX	269848.	45.813	1.69
48	106	1187	29:43	3	1.265	A BB	490632.	51.870	1.92

AR100681

00043

Water Volatile Matrix Spike/Matrix Spike Duplicate Recovery

3A

Water Volatile Matrix Spike/Matrix Spike Duplicate Recovery

Lab Name: Princeton Testing Lab Contract: WESTON, TAT

Lab Code: PTL Case No.: 9108360 SAS No.: SDG No.:

Matrix Spike - EPA Sample No.: 8360-5 (X100,000)

S4962

S4963

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC #	QC LIMITS REC
1,1-Dichloroethene	50	ND	30.1	60*	61-145
Trichloroethene	50	ND	61	122*	71-120
Benzene	50	#	# U	U	76-127
Toluene	50	3.8	54.6	102	76-125
Chlorobenzene	50	ND	48.7	97	75-130

S 4964

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD % REC #	% RPD #	QC RPD	LIMITS REC
1,1-Dichloroethene	50	30.6	61	2	14	61-145
Trichloroethene	50	61.4	123*	1	14	71-120
Benzene	50	#	U	U	11	76-127
Toluene	50	54.3	101	1	13	76-125
Chlorobenzene	50	48.8	98	1	13	75-130

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

* Values outside of QC limits

RPD: 0 out of 4 outside limits

Spike Recovery: 3 out of 8 outside limits

COMMENTS: U = Unable to quantitate due to non-homogeneous matrix

= Reproduceability of Benzene was unattainable due to non-homogeneous nature of the sample matrix.

FORM III VOA-1

1/87 Rev.

AR100682

ANALYSIS REPORT FILE SOLARIZED

Date: 8/19/1991 T1

10/20/91 12:37:00

Sample: METHOD BLANK

Column: GLASS PACKED COLUMN/METHOD 624/I.C 10/18/91

Formula: SML

Instrument: FINN

Weight: 0.000

Submitted by: PTL

Analyst: CC

Acct. No.: PTL

AMOUNT=AREA * REF AMNT/(REF AREA * RESP FACT)

Resp. fact. from Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **INT. STD. **
2	C110 1, 4-DIFLUOROBENZENE **INT. STD. **
3	C120 CHLOROBENZENE-D5 **INT. STD. **
4	CS15 1, 2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-DB **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	CO10 CHLOROMETHANE **
8	C015 BROMOMETHANE
9	C020 VINYL CHLORIDE *
10	C025 CHLOROETHANE
11	C030 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1, 1-DICHLOROETHENE *
17	C050 1, 1-DICHLOROETHANE **
18	C055 TRANS-1, 2-DICHLOROETHENE
19	C000 TRICHLOROFUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1, 2-DICHLOROETHANE
22	C110 2-BUTANONE
23	C115 1, 1, 1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1, 2-DICHLOROPROPANE *
28	C145 TRANS-1, 3-DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1, 1, 2-TRICHLOROETHANE
32	C165 BENZENE
33	C140 CIS-1, 3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C240 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1, 1, 2, 2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M-XYLENE
45	C250 1, 3-DICHLOROBENZENE
46	C254 1, 4-DICHLOROBENZENE
47	C255 1, 2-DICHLOROBENZENE

AR100683

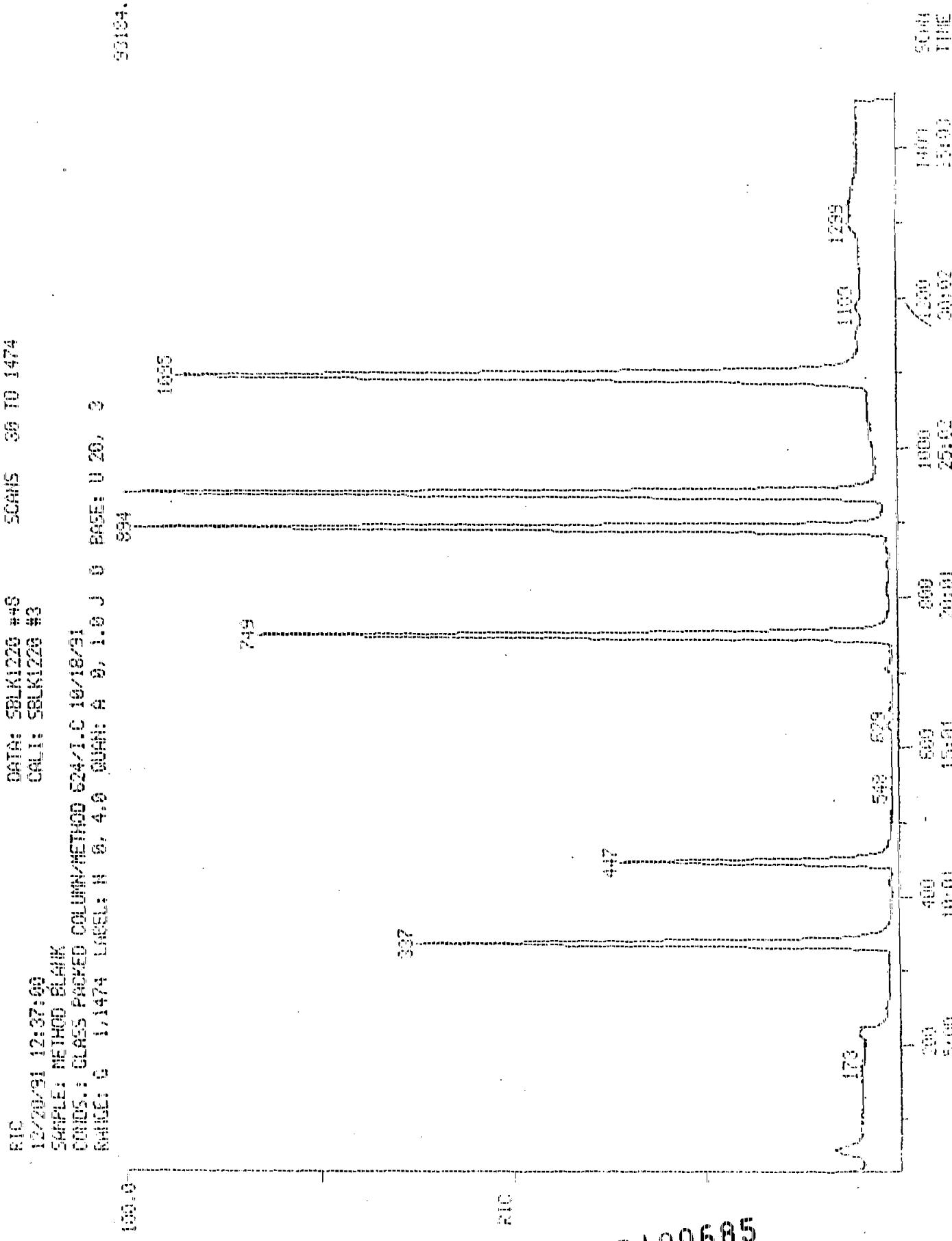
00051

No. Blanks
43 0200 C4P XYLYENE

No.	m/z	Scan	Time	Ref	RRT	Meth	Area(Hight)	Attenuat	% Tot
1	49	337	8:26	1	1.000	A BB	123077.	50.000 NG	16.37
2	114	737	13:45	2	1.000	A BB	267686.	50.000 NG	16.37
3	117	939	23:30	3	1.000	A BB	251449.	50.000 NG	16.37
4	65	447	11:11	1	1.326	A BB	86906.	40.544 NG	14.26
5	98	874	22:23	3	0.952	A BB	275915.	51.432 NG	15.85
6	95	1026	27:26	3	1.167	A BB	203909.	51.652 NG	16.91
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	49	217	8:26	1	0.644	A BB	4179.	2.556 NG	✓ 0.84
12	NOT FOUND								
13	43	251	6:17	1	0.745	A BB	1454.	4.379 NG	✓ 1.43
14	NOT FOUND								
15	NOT FOUND								
16	NOT FOUND								
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	43	536	13:25	2	0.716	A BB	79.	0.080 NG	0.02
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	95	630	15:45	2	0.841	A BB	1075.	0.528 NG	0.18
30	NOT FOUND								
31	NOT FOUND								
32	NOT FOUND								
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	NOT FOUND								
37	NOT FOUND								
38	43	810	20:16	3	0.863	A BB	905.	0.254 NG	0.26
39	NOT FOUND								
40	91	901	22:03	3	0.960	A BB	1470.	0.266 NG	0.08
41	NOT FOUND								
42	106	1016	25:26	3	1.082	A BB	419.	0.194 NG	0.06
43	NOT FOUND								
44	NOT FOUND								
45	NOT FOUND								
46	NOT FOUND								
47	NOT FOUND								
48	NOT FOUND								

AR100684

00051



ANALYST: FULTONATIC

DIAGNOSTIC REPORT

JUL 21 1984 8:47:02

DATA FILE: GOLK12.D

FILTER SCAN PARAMETERS

MAX NUMBER TICS: 15
TBL-TABLE ENTRIES: 528
SCAN TOLERANCE : 2
MIN RIC HT. DDX: 10
FIRST SCAN : 1
LAST SCAN : 1600
TIC THRESHOLD : 600

NOTIFICATION LIBRARIES & LISTS

TIC LIBRARY: LIBRARYLIB
NBS SEARCH PROC: BERLIB
PEAK FINDER PROC: VCOMC
TCA TIC LL : LS
FILE NAME LIST : TCAREF2

TARGET COMPOUND ANALYSIS:

TARGETS (QUAN LIST)	IS PEAKS	TOTAL TARGET PEAKS
10	9	19

FILTER PROCESSING:

REJECT PEAKS							
TOTAL PEAKS	< 1ST SCAN	> LAST SCAN	< MIN RIC HT	< SCAN TOL	> MAX # PEAKS	TOTAL REJECTS	TOTAL TICS
6	0	0	0	6	0	6	0

NO UNKNOWN PEAKS TO BE IDENTIFIED.

ARI00686

00054

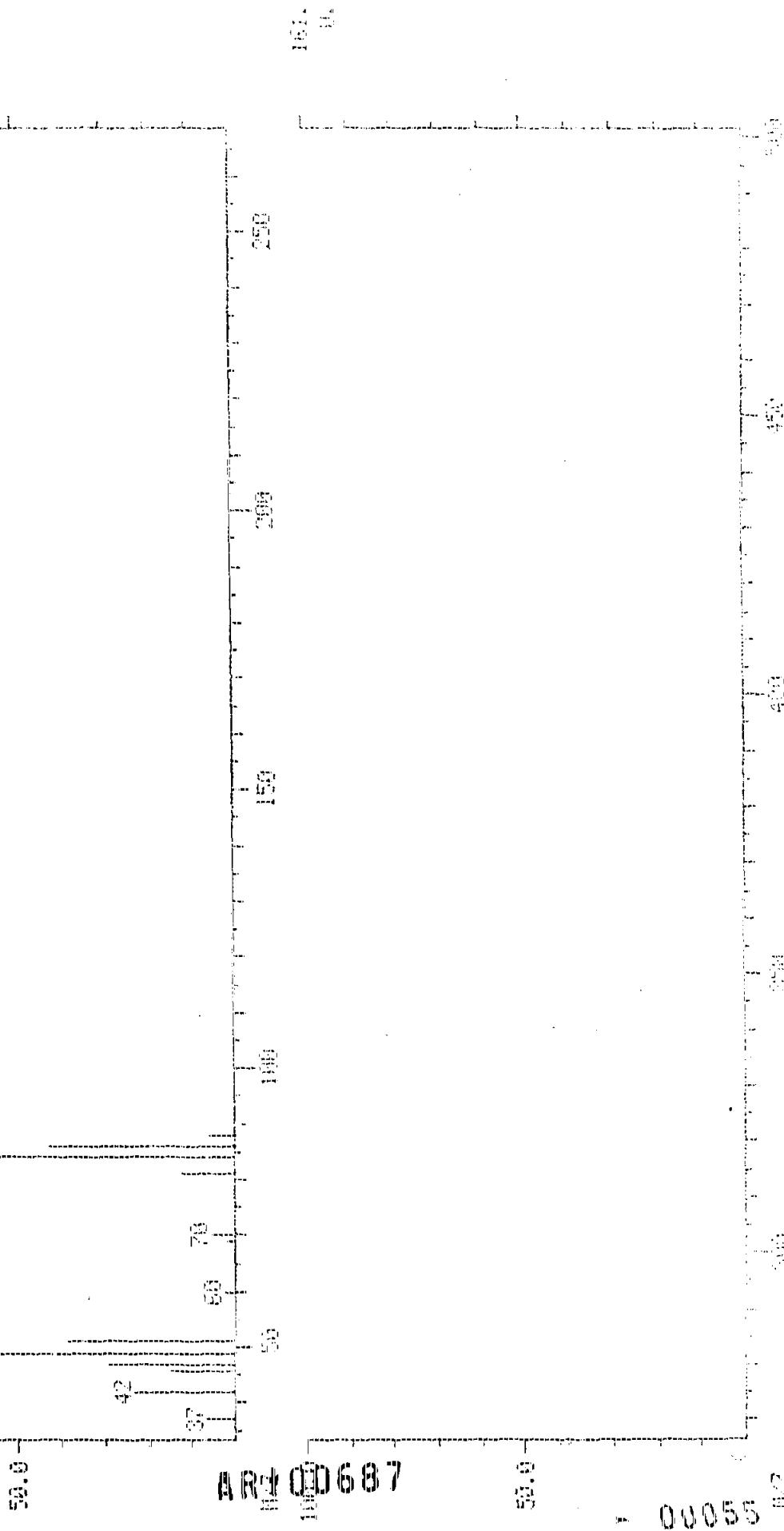
MASS SPECTRUM
12/26/91 12:37:69 + 5:26
SAMPLE: METHANOL BLANK
COLLECT: GLASS PACKED COLUMN/METHYL B24/1, C 10/16/91
TEMP: 54.000, C
EQUIPPED (5) 100 2H 9T)

DATA: SSI K1229 #217 BASE M/Z: 49
CAL: SSI K1229 #3 RIG: 553,

100.0

SI

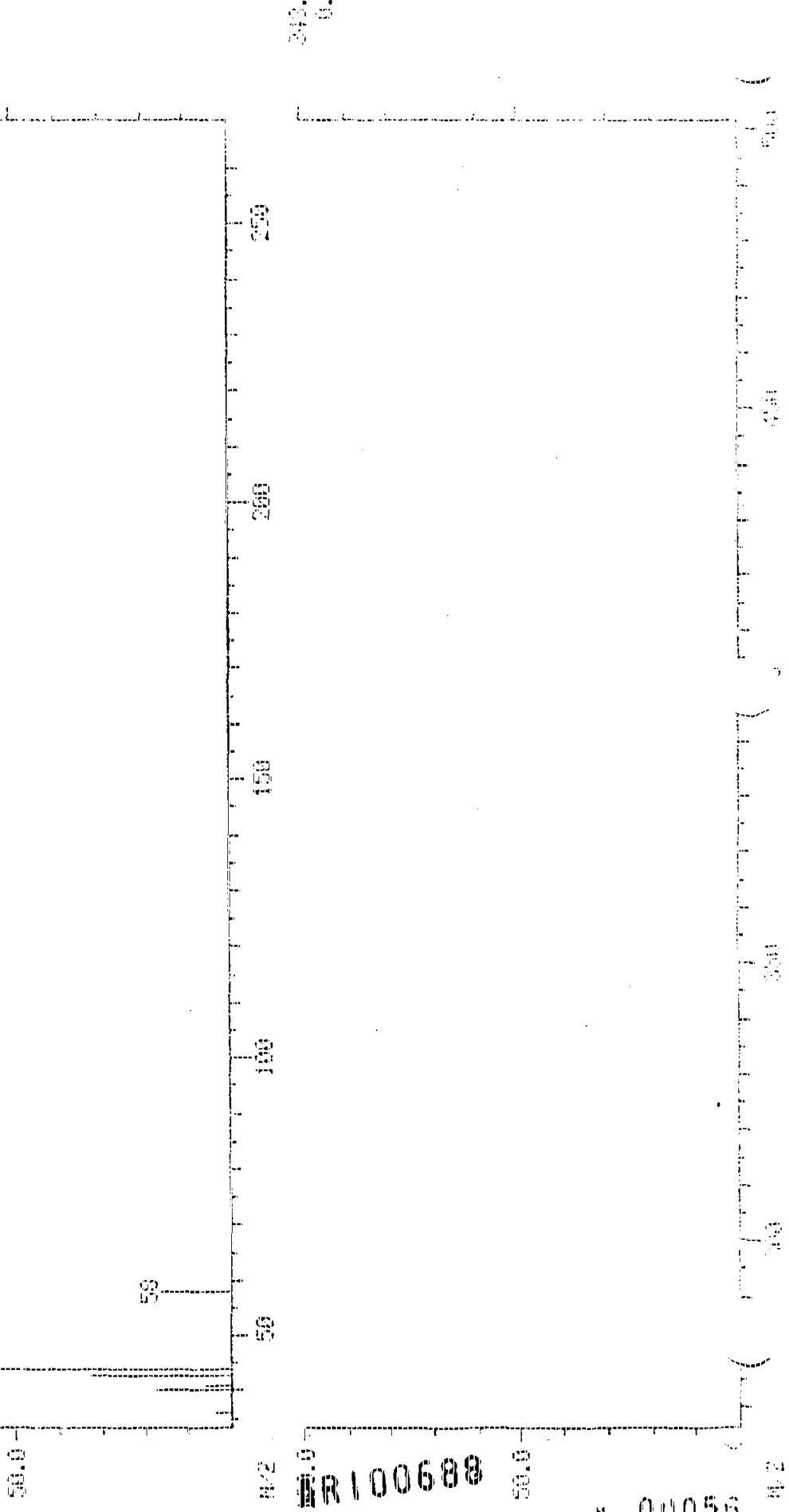
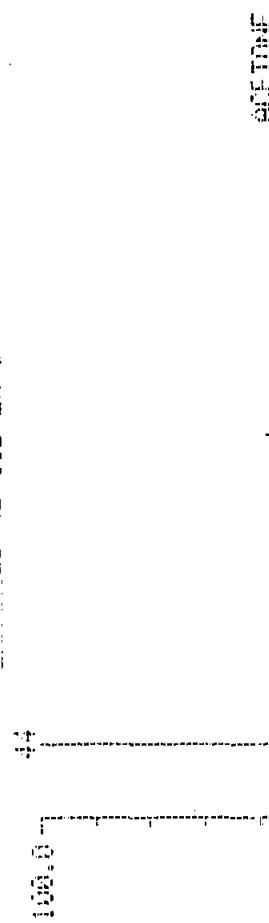
NEUTRAL CHARGE



MASS SPECTRUM
12/20/91 12:37:00 + 5:17
SAMPLE: METHYL BIS(ACRYLATE)
CONC.: 0.005% (VOL/VOL) 524/1.0 10/18/91
TEMP: 71 DEG. C
EQUIPMENT (S 100 20 97)

DATA: SBK1229 #251
CALL: SBK1229 #3

BASE: #2: 44
RT0: 50.4.



Identification Report File: SOLKID

Date: 10/10/82

Time: 12:15:00

Sample: METHYL CHLORIDE

Location: PLATE PACIFIC OCEAN, LAT 30° S 62° E

Instrument: FID/NI

Analytical Method: GC

Dilution: 1:200

Purge: N2, PTL

AMOUNT=AREA * REF AMNT/REF AREA * RESP FACT
Resp. Fac. From Library Entry

No	Name
1	C101 BROMOCHLOROMETHANE **INT. STD. **
2	C110 1, 4-DIFLUOROBENZENE **INT. STD. **
3	C120 CHLOROBENZENE-D6 **INT. STD. **
4	CS15 1, 2-DICHLOROETHANE-D4 **S. STD. **
5	CS05 TOLUENE-D8 **S. STD. **
6	CS10 4-BROMOFLUOROBENZENE **S. STD. **
7	CO10 CHLOROMETHANE **
8	CO15 BROMOMETHANE
9	CO20 VINYL CHLORIDE *
10	CO25 CHLOROETHANE
11	CO30 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C035 ACETONE
14	C252 ACRYLONITRILE
15	C040 CARBON DISULFIDE
16	C045 1, 1-DICHLOROETHENE *
17	C050 1, 1-DICHLOROETHANE **
18	C055 TRANS-1, 2-DICHLOROETHENE
19	C000 TRICHLOROFUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1, 2-DICHLOROETHANE
22	G110 2-BUTANONE
23	C115 1, 1, 1-TRICHLOROETHANE
24	C120 CARBON TETRACHLORIDE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1, 2-DICHLOROPROPANE *
28	C145 TRANS-1, 3-DICHLOROPROPENE
29	C150 TRICHLOROETHENE
30	C155 DIBROMOCHLOROMETHANE
31	C160 1, 1, 2-TRICHLOROETHANE
32	C165 BENZENE
33	C140 C15-1, 3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1, 1, 2, 2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C250 M-XYLENE
45	C250 1, 3-DICHLOROBENZENE
46	C254 1, 4-DICHLOROBENZENE
47	C255 1, 2-DICHLOROBENZENE

AR100689

00057

11 Notes
12 Gluco C13 Valence

Line	Row#	Scan	Time	Rat#	HPT	Net%	Area (Integration)	Average	NTot
1	49	2007	5.22	2	0.600	A 60	1984.07	50.010 NG	15.42
2	104	749	13:45	2	0.610	A 60	5637.2	50.010 NG	15.42
3	117	762	21:30	2	0.600	A 60	5013.07	50.000 NG	15.42
4	63	467	11:11	1	0.226	A 60	1261.73	50.000 NG	15.22
5	93	374	23:20	3	0.752	A 60	5478.97	45.000 NG	15.38
6	95	1035	27.26	3	1.167	A 60	3808.00	50.010 NG	16.09
7	50	49	1:14	1	0.145	A 60	2179.	0.500 NG	0.16
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	49	216	5.64	1	0.641	A 60	4747.	4.519 NG	0.47
12	NOT FOUND								
13	43	257	6.66	1	0.763	A 60	7797.	12.453 NG	3.64
14	NOT FOUND								
15	NOT FOUND								
16	NOT FOUND								
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	97	498	12:28	2	0.665	A 60	449.	0.099 NG	0.03
24	NOT FOUND								
25	NOT FOUND								
26	NOT FOUND								
27	NOT FOUND								
28	NOT FOUND								
29	95	631	15:48	2	0.842	A 60	2137.	0.509 NG	0.16
30	NOT FOUND								
31	NOT FOUND								
32	79	652	16:19	2	0.870	A 60	3259.	0.293 NG	0.09
33	NOT FOUND								
34	NOT FOUND								
35	NOT FOUND								
36	166	844	21:08	3	0.899	A 60	2361.	0.395 NG	0.12
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	91	901	22:33	3	0.950	A 60	5209.	0.420 NG	0.13
41	112	943	23:35	3	1.004	A 60	4375.	0.427 NG	0.13
42	109	1015	23:36	3	1.062	A 60	2253.	0.510 NG	0.16
43	104	1142	23:44	3	1.223	A 60	4077.	0.429 NG	0.15
44	105	1176	23:56	3	1.231	A 60	2915.	0.501 NG	0.19
45	145	1283	32:29	3	1.382	A 60	2305.	0.489 NG	0.15
46	146	1329	33:11	3	1.412	A 60	349.	0.691 NG	0.08
47	146	1343	33:44	3	1.426	A 60	1536.	0.500 NG	0.09
48	106	1135	29:41	3	1.263	A 60	4507.	0.429 NG	0.14

AR100690

00055

RIC
01/20/2022 12:15:00
SAMPLE: METHOD BLANK

COLLECT: GLASS PACKED COLUMN/METHOD 624

POLAR: G 1,1474 LABEL: H 6, 4.9 DETEC: A 6, 1-3 J 9 845E: U 29, 3

162272.

DATA: SELK12 #48
CALC: SELK12 #3

SCALES 30 70 1474
30445 30 70 1474
30445 30 70 1474

100.0

75.0

50.0

25.0

0.0

385

437

547

631

844

1018

1153

1220

1418

1560

1630

1700

1000 /
2000 /
3000 /
4000 /
5000 /
TIME

1000
2000
3000
4000
5000

1000
2000
3000
4000
5000

1000
2000
3000
4000
5000

1000
2000
3000
4000
5000

1000
2000
3000
4000
5000

1000
2000
3000
4000
5000

PPR00691

PPR00691

DATA FILE: SELK12 #215
TARGET COMPOUND COMPARISON
COMPOUND: 0335 METHYLENE CHLORIDE

STANDARD FILE: SU0012R #217
CAL I: SELK12 #3

PUB DATA: SELK12 #215
01/02/92 12:15

BASE M/Z: 44 RIC: 5120.

2552.

20.0

10.0

STANDARD FILE: SELK12 #215
01/02/92 11:09

BASE M/Z: 44 RIC: 5120.

2552.

10.0

5.0

1.0

0.5

0.2

0.1

0.05

0.02

0.01

0.005

0.002

0.001

0.0005

0.0002

0.0001

-0.0002

-0.0005

00692

BASE M/Z: 49 RIC: 27456.
CHAMACED (5 158 21:07)

2552.

100.0

80.0

60.0

40.0

20.0

0.0

-20.0

-40.0

-60.0

-80.0

000

DATA FILE: SELK12 #257
TARGET COMPiling COMPUTER 1500
CPU/WORD: 0.035 SEC/TIME

STANDARD FILE: SV012R #250
CALL: SELK12 #2

DATA FILE: SELK12 #257
01/02/92 10:15

DATA FILE: SELK12 #257
01/02/92 10:15

29.17

14.56

CONTINUED FILE: SELK12 #257

DATA FILE: SELK12 #257
01/02/92 10:15

14.57

7.8

4R - 0 00
166 93

STANDARD FILE: SV012R #250
01/02/92 10:15

DATA FILE: SELK12 #257
01/02/92 10:15

7.8

0.6

0.2

166.0

50.0

0.6

-50.0

-166.0

0.0065

Laboratory Report File 3470

Date: 3470-71
12-10-68 15 07.30

Sample: 6360-1

Column: GLASS PACKED COLUMN METHOD 624/1 C 100/10%
Furnace: 200°, RF 4000V/BML Instrument: FID

Submitted by: WESTON Analyst: JC

Weight: 0.400

Act No.: 6360

CHLORINATED & RELATED AREA & REFL FACTS

Keep: Acc Print Library, Entry

No.	Name
1	C101 SPONCHILODENE THANE ** INT STD. **
2	C110 1,4-DIFLUORODENEINE ** INT. STD. **
3	C120 CHLOROBENZENE-DB ** INT. STD. **
4	C130 1,2-DICHLOROETHANE-04 ** STD. **
5	C135 TOLUENE-DB **G. STD. **
6	C140 4-BROMOFLUOROBENZENE **B. STD. **
7	C150 CHLOROMETHANE **
8	C160 BROMOMETHANE
9	C170 VINYL CHLORIDE
10	C175 CHLOROETHANE
11	C180 METHYLENE CHLORIDE
12	C185 ACROLEIN
13	C190 ACETONE
14	C200 1,1,2,2-TETRACHLOROETHANE
15	C240 CARBON DISULFIDE
16	C245 1,1-DICHLOROETHENE *
17	C250 1,1-DICHLOROETHANE **
18	C255 TRANS-1,2-DICHLOROETHENE
19	C260 TRICHLOROFLUOROMETHANE
20	C265 CHLOROFORM *
21	C285 1,2-DICHLOROETHANE
22	C290 1,1,2,2-TETRACHLOROETHANE
23	C310 1,1,2,2-TETRACHLOROETHANE
24	C320 1,1,2,2-TETRACHLOROETHANE
25	C125 VINYL ACETATE
26	C130 BROMO DICHLOROMETHANE
27	C140 1,2-DICHLOROPROPANE
28	C145 TRANS-1,2-DICHLOROPROPENE
29	C150 PROPENE
30	C155 1,1,2,2-TETRACHLOROETHANE
31	C160 1,1,2,2-TETRACHLOROETHANE
32	C165 TETZENE
33	C170 CIS-1,3-DICHLOROPROPENE
34	C175 2-CHLOROETHYL VINYL ETHER
35	C180 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C225 4-METHYL 2-PENTANONE
39	C226 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C245 CHLOROBENZENE **
42	C250 ETHYL CHLORIDE *
43	C260 CYCLOPENTANE
44	C270 CYCLOPENTENE
45	C280 CYCLOPENTANE
46	C290 CYCLOPENTENE

AP100694

0006

10063
10064 C.R.P. ANALYST

#	Ref	Sample	Time	Ref#	PRT	Match	Area (High)	Acq. Date	Acq. Time
1	49	642	8:34	1	1.000	A 28	118028	50.000 NG	1.32
2	114	751	18:48	2	1.000	A 28	668886	50.000 NG	1.32
3	117	758	20:29	3	1.000	A 28	377857	50.000 NG	1.32
4	66	465	11:08	4	1.001	A 28	104956	50.112 NO	1.94
5	98	682	22:20	5	0.951	A 28	361841	47.912 NO	1.73
6	98	1054	27:00	6	1.164	A 28	290521	43.487 NO	1.77
7	50	84	11:17	7	0.149	A 28	2288	43.140 NO	0.06

8 NOT FOUND

9 NOT FOUND

10 NOT FOUND

11 49 642 8:34 1 0.642 A 28

Acq. Date	Acq. Time
50.000 NG	1.32
50.000 NG	1.32
50.000 NG	1.32
50.112 NO	1.94
47.912 NO	1.73
43.487 NO	1.77
43.140 NO	0.06

12 NOT FOUND

13 NOT FOUND

14 66 465 11:08 2 1.001 A 28

50.112 NO ✓ 0.21

15 NOT FOUND

16 NOT FOUND

17 NOT FOUND

18 NOT FOUND

19 NOT FOUND

20 NOT FOUND

21 NOT FOUND

22 50 84 11:17 2 0.149 A 28

44100000
50.140 NO ✓ 0.44

23 NOT FOUND

24 NOT FOUND

25 NOT FOUND

26 NOT FOUND

27 NOT FOUND

28 NOT FOUND

29 NOT FOUND

30 NOT FOUND

31 NOT FOUND

32 78 650 10:16 2 0.646 A 28

50.112 NO ✓ 0.08

33 NOT FOUND

34 NOT FOUND

35 NOT FOUND

36 NOT FOUND

37 NOT FOUND

38 NOT FOUND

39 NOT FOUND

40 49 675 22:24 2 0.768 A 28

50.112 NO ✓ 0.16

41 NOT FOUND

42 106 1016 22:24 2 1.092 A 28

50.112 NO ✓ 0.38

43 NOT FOUND

44 106 1156 22:26 2 1.002 A 28

50.112 NO ✓ 0.94

45 NOT FOUND

46 NOT FOUND

47 NOT FOUND

48 106 1156 22:41 2 1.264 A 28

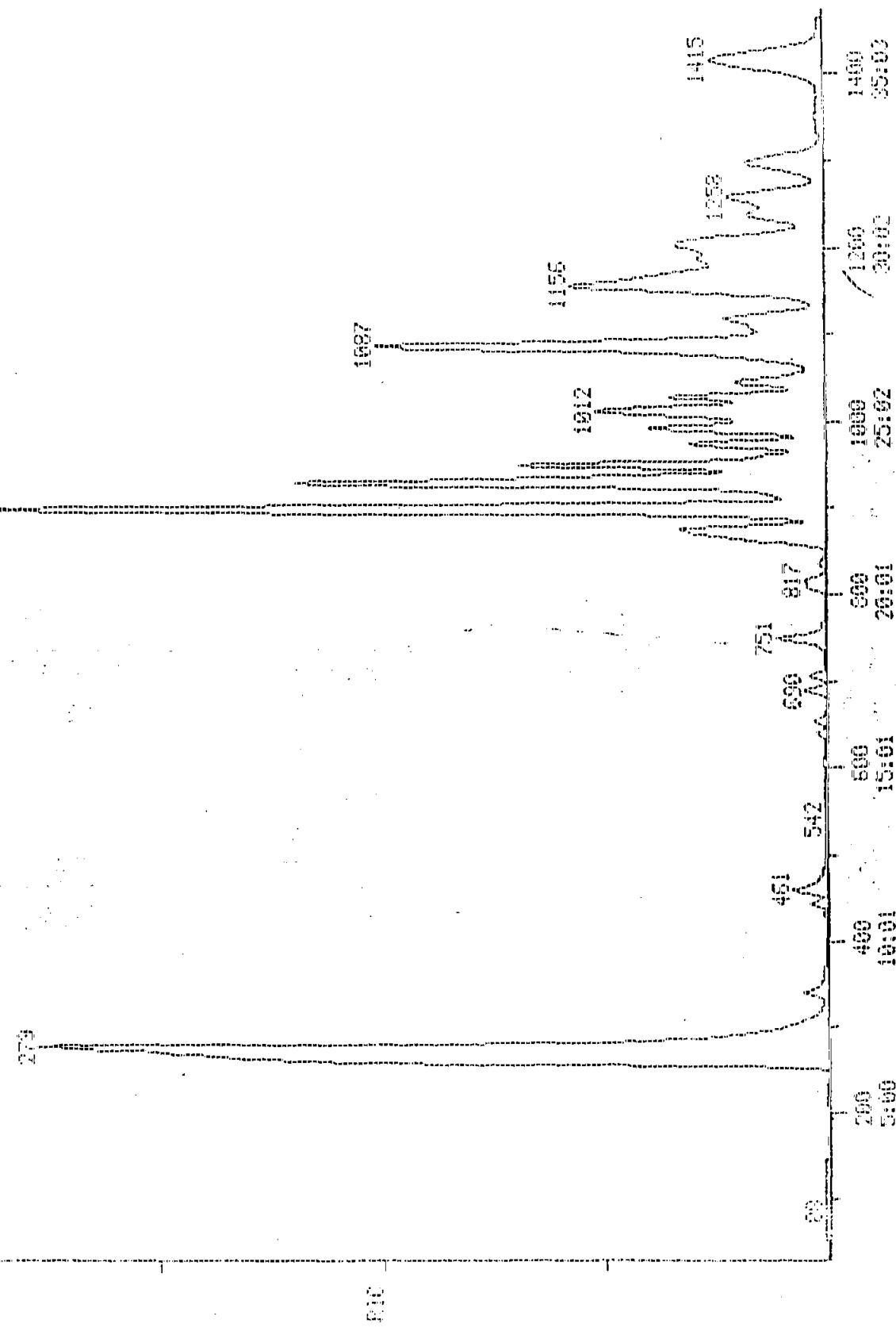
50.107 NO ✓ 0.29

PR100695
00063

RIE

12/29/91 15:07:00
SAMPLE: S369-1/
COLUMN: GLASS PACKED COLUMN/METHOD 624/1.C 10/18/91
PAGE: C 1,1474 LABEL: N , 0, 4.0 0.0000 0 0.1,5 0 0.0000 0 0.20, 0
2291470.

DATA: 54900 #46
SCANS: 39 TO 1474



AR 100696

000000000000

MS5 SPECTRUM
12/26/91 15:07:06 + 1:15
SAMPLE: 8366-1
09405.: GLASS PLATED COLUMN/METHOD 524/1.C 12/19/91
TEMP: 45 SEC. C

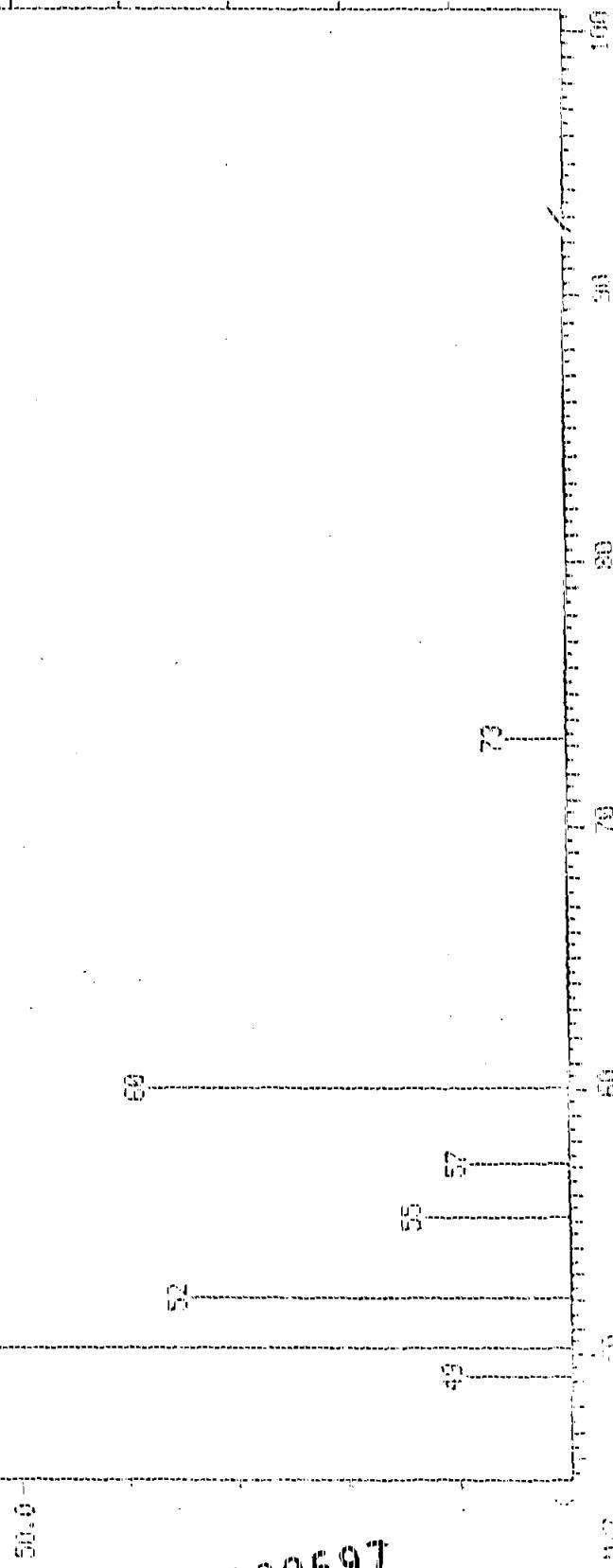
50

100.0

273.
9.

GR/EDGE
=9.6584

CHLOROMETHANE



DATA FILE: 54200 #52
TARGET COMPOUND COMPARISON
COMPOUND: C9H10 CHLOROMETHANE
PAN DATE: 54200 #52
12/26/91 15:07

STANDARD FILE: 540122R #48
CALL: 54200 #3

BASE W/Z: 44

RIC: 2672.

300.0
200.0
100.0
0.0

ENHANCED DATA: 54200 #52

STANDARD FILE: 540122R #48
12/26/91 15:00

BASE W/Z: 44

RIC: 2546.

150.0
100.0
50.0
0.0

STANDARD FILE: 540122R #48
12/26/91 15:00

BASE W/Z: 39

ENHANCED (S 158 26 OT)

100.0
50.0
0.0

AR 100698

00066

DATA FILE: S4306 #226 STRANGL. : n/a: SU01220R #217
TARGET COMPOUND COMPARISON
COMPODE: 0326 METHYLENE CHLORIDE

FILE DATE: 12/28/91 15:47
S4306 #226

BASE M/Z: 44 RIC: 5704.



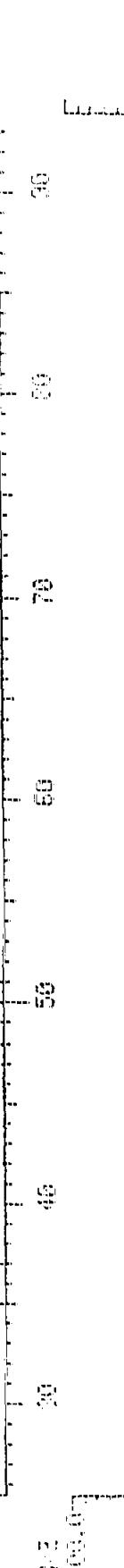
ENHANCED DATA: S4306 #226

BASE M/Z: 43 RIC: 5704.



STRANGL. FILE: SU01220R #217
12/28/91 15:46

BASE M/Z: 45 RIC: 5704.



PR 100699

-100.0

0.0

00067

DATA FILE: S4900 #451
TARGET COMPOUND COMPARISON!
COMPOUND: C116 2-EUROHEX
FROM: 90101: S4900 #451
12/26/91 15:07

STANDARD FILE: S001228R #466
CBL1: S4900 #2

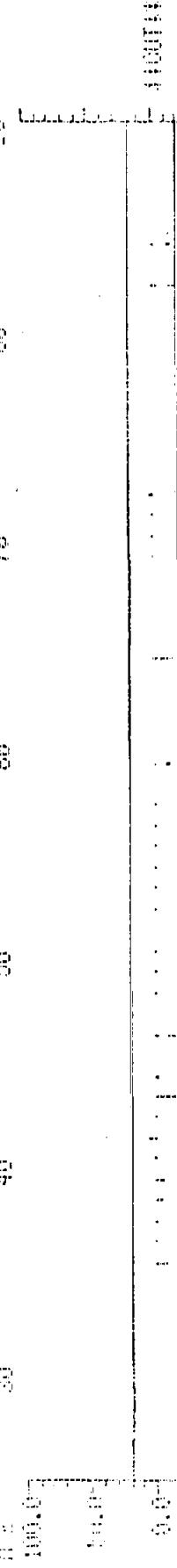
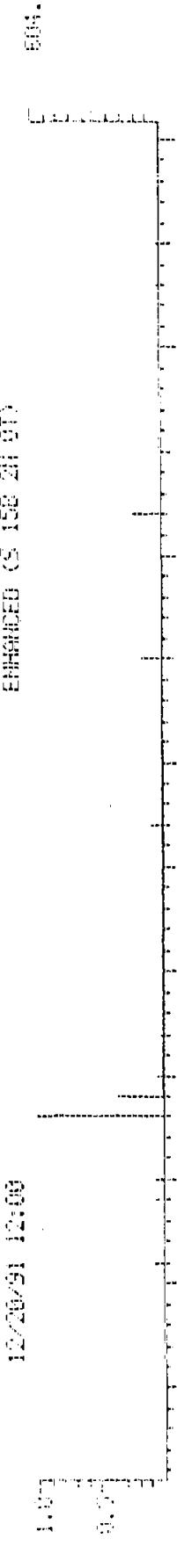
BASE M/Z: 42

BASE M/Z: 43

100.0
80.0
60.0
40.0
20.0
0.0

STANDARD FILE: S001228R #466
12/26/91 12:00

STANDARD FILE: S001228R #466
12/26/91 12:00



AR100700

00065

DATA FILE: 54906 #893
TARGET COMPOUND: C236 TOLUENE
COMPOUND: C236 TOLUENE *

STANDARD NAME: SUDANONE #901
CHL: 54906 #8

FILE DATE: 5/29/91 15:07
FILE NUMBER: 54906 #893
PAGE N/2: 91
PICS: 2255670.

SCANNED.

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BR100701

00069

DATA FILE: S4300 #1015
TARGET COMPOUND: COPPER(II)
COMPOUND: C240 ETHYL BENZENE *

STANDARD FILE: S401208 #1015
STANDARD: S4300 #10

PPM: 6074: 34300 #1015
12/20/91 15:07

BASE W/Z: 91

100.0

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STANDARD FILE: S401208 #1015
STANDARD: S4300 #10

PPM: 6074: 34300 #1015
12/20/91 15:07

BASE W/Z: 91

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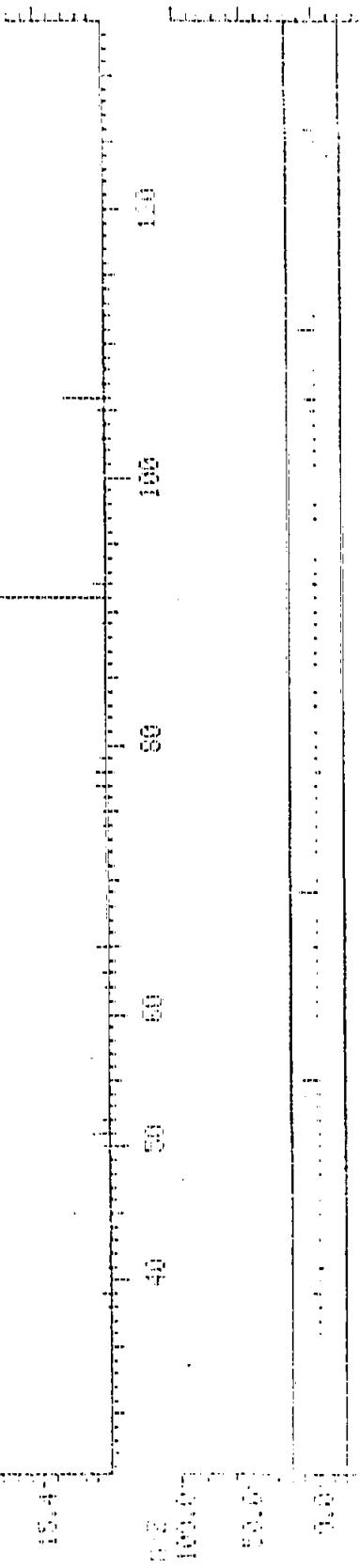
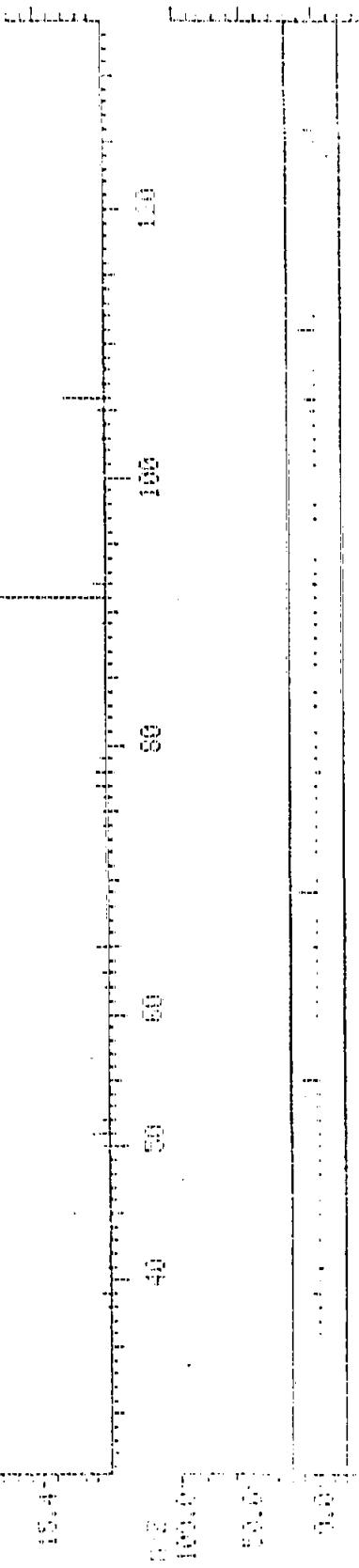
100.0

100.0

STANDARD FILE: S401208 #1015
12/20/91 15:07

BASE W/Z: 91
EQUATED CS 158 21 OT?

32500,



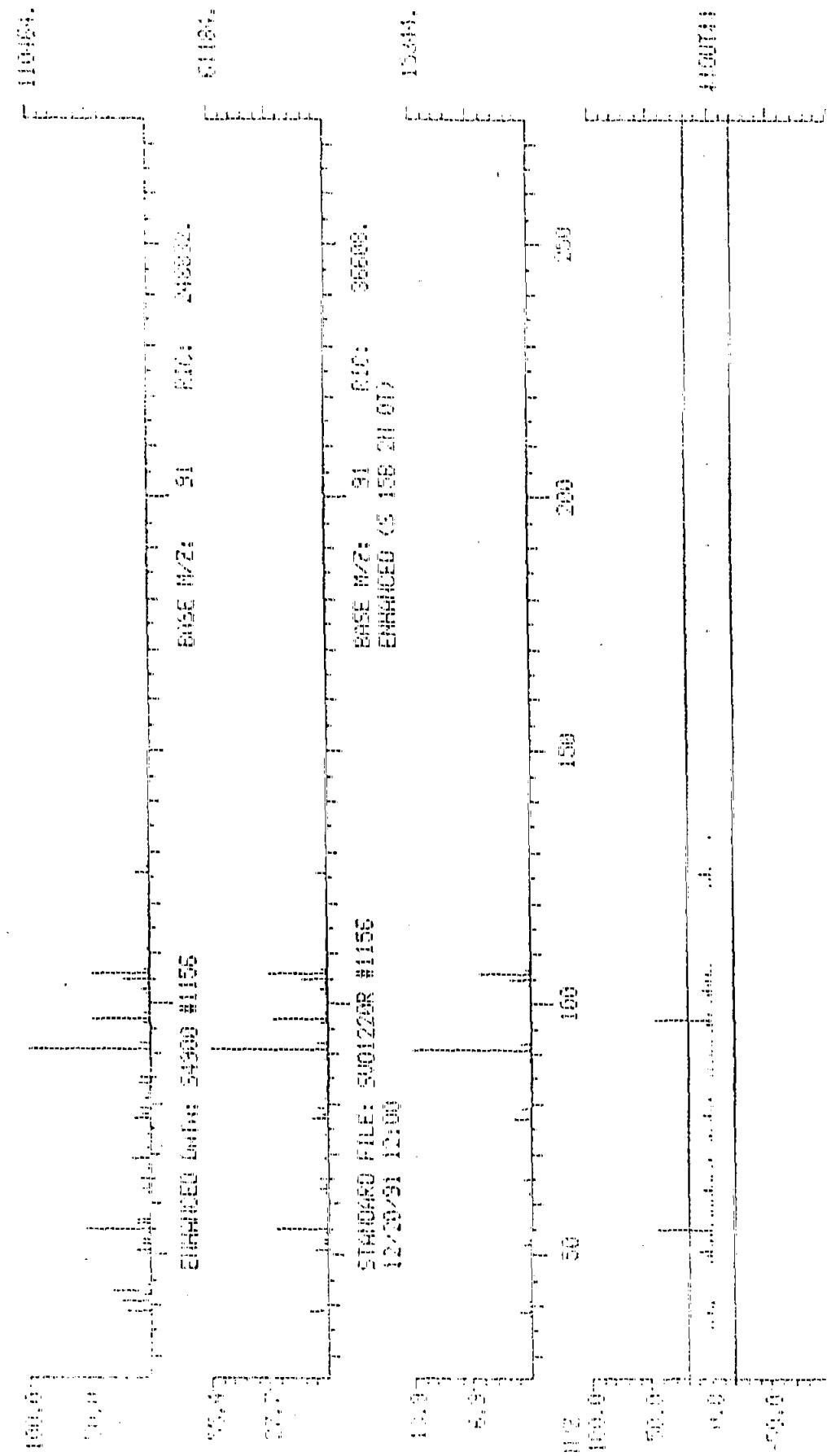
AR100702

00079

901A FILE: 54990 #1156 SOURCE FILE: SH012208 #1156
TARGET COMPOUND COMPARISON
COMPOUND: C250 N-NYLENE

REC DATE: 5/29/91 15:07 BASE M/Z: 91 RIC: 545120

BASE M/Z: 91 RIC: 545120



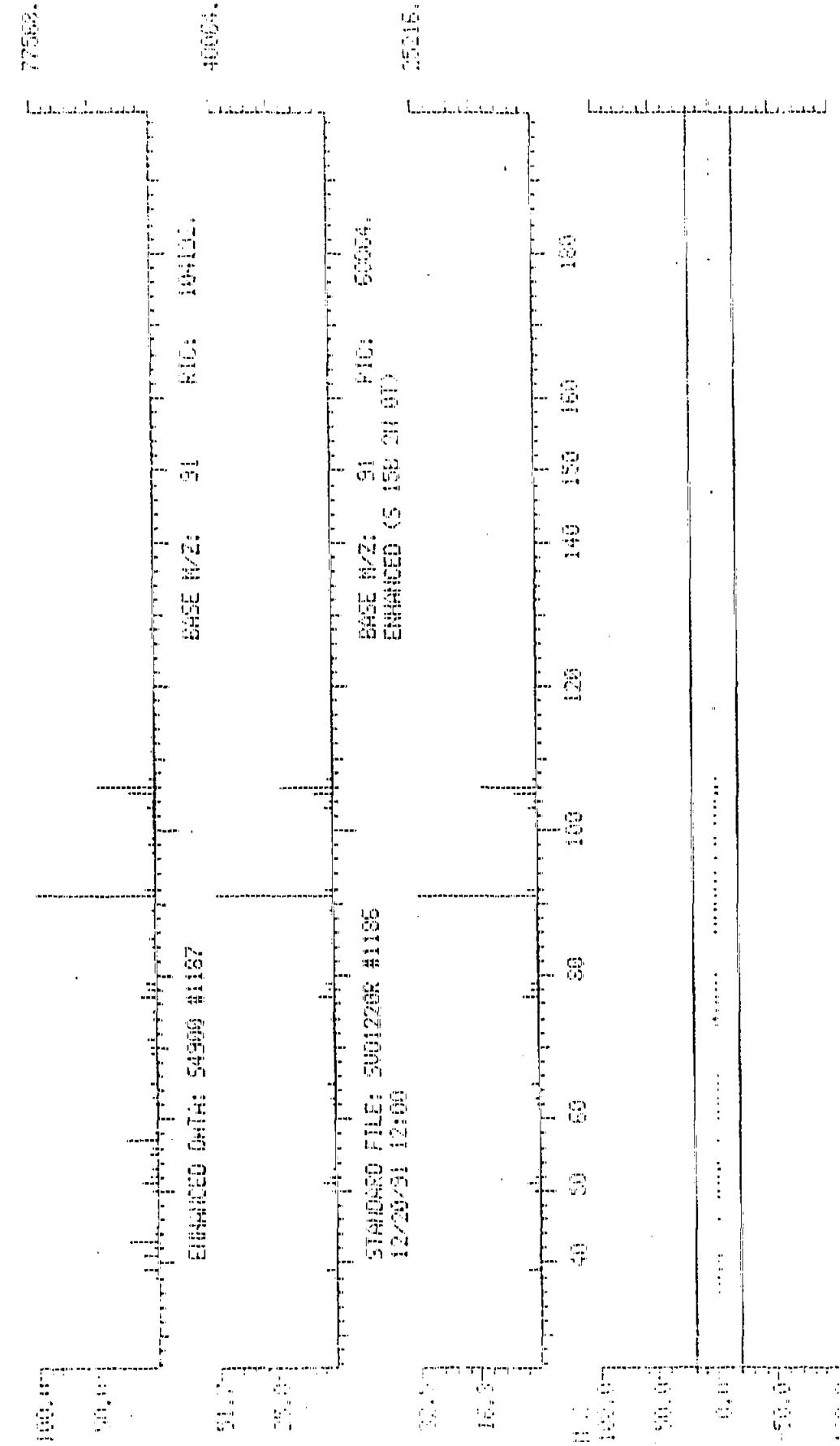
AB100703

00071

DATA FILE: S4999 #1187
TARGET COMPOUND COMPARISON
COMPOUND: C250 8AP XYLENE
REF. DATA: S4999 #1187
12/20/91 15:07

STANDARD FILE: S001220R #1185
CALI: S4999 #2

BASE M/Z: 91 RIC: 317652.



MR100704

00072

Identification Report File: G4PC

Date: 8402.11

12000. PI 16:36:00 *4040*

Devicer: GC602

Column: GLASS PACKED COLUMN/METHOD #24/I.C. 10/12/81

For user: 100UL DEDOX/5ML Instrument: FINN

Published by: WESTON Analyst: CC

Report Date: 6.26.00

Page No.: 1/16

CHLORINEAREA = REF AMNT/(REF AREA * REEF FACT)

Reef. fact. from library entry

#	Name
1	C101 BROMOCHLOROMETHANE **INT. STD. **
2	C110 1,4-DIFLUOROBENZENE **INT. STD. **
3	C120 CHLOROBENZENE-08 **INT. STD. **
4	C135 1,2-DICHLOROETHANE-04 **S. STD. **
5	C205 TOLUENE-08 **S. STD. **
6	C210 4-BROMOFLUOROBENZENE **S. STD. **
7	C210 CHLOROMETHANE **
8	C215 BROMOMETHANE
9	C220 VINYL CHLORIDE *
10	C225 CHLOROETHANE
11	C230 METHYLENE CHLORIDE
12	C251 ACROLIN
13	C255 ACETONE
14	C259 ACRYLONITRILE
15	C260 CHLORO BISULFIDE
16	C265 1,1-DICHLOROETHENE *
17	C270 1,1-DICHLOROETHANE **
18	C275 TRANS-1,2-DICHLOROETHENE
19	C280 TRICHLOROPROPROMETHANE
20	C280 CHLOROFORM *
21	C285 1,2-DICHLOROETHANE
22	C290 2-BUTANONE
23	C290 CARBON TETRACHLORIDE
25	C295 VINYL ACETATE
26	C300 BROMO DICHLOROMETHANE
27	C340 1,2-DICHLOROPROPANE *
28	C345 TRANS-1,3-DICHLOROPROPENE
29	C350 TRICHLOROETHENE
30	C365 DIBROMOCHLOROMETHANE
31	C365 1,1,2-TRICHLOROETHANE
32	C365 BENZENE
33	C370 CIS-1,3-DICHLOROPROPENE
34	C375 2-CHLOROETHYL VINYL ETHER
35	C380 BROMOFORM **
36	C220 TETRACHLOROETHENE
37	C210 2-HEXANONE
38	C205 4-METHYL 2-PENTANONE
39	C225 1,1,2,2-TETRACHLOROETHANE **
40	C230 TOLUENE *
41	C235 CHLOROBENZENE **
42	C240 ETHYL BENZENE *
43	C245 STYRENE
44	C245 PROPENE
45	C250 PROPENE
46	C250 PROPENE

AR100705

• 00073

Data: C6B XYLENE

	Run	Span	Time	Ref	RPT	Match	Area (High)	Accuracy	%Tot
1	109	837	6:26	1	1.000	A BB	102664.	50.000 NG	2.00
2	114	749	18:46	2	1.000	A BB	369524.	50.000 NG	2.00
3	117	938	20:29	3	1.000	A BB	353425.	50.000 NG	2.00
4	65	444	11:07	1	1.318	A BB	102606	59.152 NG	2.40
5	98	893	22:21	3	0.952	A BB	395724.	52.486 NG	2.13
6	95	1094	27:23	3	1.166	A BB	275531.	49.178 NG	1.99
7	NOT FOUND								
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	109	816	6:24	1	0.641	A BB	102114	50.512 NG	0.34
12	NOT FOUND								
13	43	858	6:27	1	0.766	A BB	16166	57.488 NG	2.00
14	NOT FOUND								
15	NOT FOUND								
16	NOT FOUND								
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	NOT FOUND								
26	106	846	21:11	3	0.902	A BB	1002.	50.000 NG	0.01
27	NOT FOUND								
28	NOT FOUND								
29	NOT FOUND								
30	NOT FOUND								
31	NOT FOUND								
32	NOT FOUND								
33	NOT FOUND								
34	NOT FOUND								
35	91	901	22:33	3	0.961	A BB	453232.	55.473 NG	2.25
36	NOT FOUND								
37	106	1014	25:23	3	1.081	A BB	2141500.	705.750 NG	28.59
38	NOT FOUND								
39	106	1154	28:53	3	1.230	A BB	3531540.	997.502 NG	40.44
40	NOT FOUND								
41	NOT FOUND								
42	NOT FOUND								
43	106	1185	29:40	3	1.263	A BB	2277840.	201.055 NG	10.42

AR100706

00074

PT
12-20-81 16:38:00
SAMPLE: 3950 ~~✓~~ CC
COMPO: 0.005 PERCENT COLIFORM/NETTOD 52471.C 10/18/81
PENIT: 6 1,474 LABEL: N 0, 4.0 QUANT: A 0, 1.0 J 0 BASE: 0 20, 3
100.0

00100 34992 448

CAL1: 34992 43

5247460.

100

BR100707

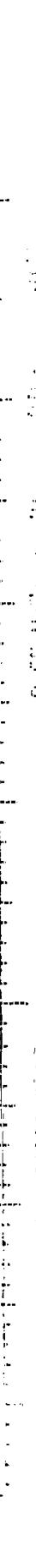
00075

NAME PLATE: 54902 #216
TARGET COMPOSITION COMPARISON
COMPTAHO: CO₂ METHYLENE CHLORIDE
PRO DATE: 12/26/91 16:36

STANDARD FILE: 54902 #216

CAL1: 54902 #2

SAGE M/Z: 49 RIC: 50.03.



AR100708

00076

DATA FILE: 542200 4200
DISSET COMMUNICATON
COMPUTER USES ALPHABET

DATA FILE: 542200 1233
DISSET COMMUNICATON
12/20/91 1513

DATA FILE: 542200 4200
DISSET COMMUNICATON
COMPUTER USES ALPHABET

DATA FILE: 542200 4200
DISSET COMMUNICATON

DATA FILE: 542200 4200
DISSET COMMUNICATON
COMPUTER USES ALPHABET

DATA

FILE

4200

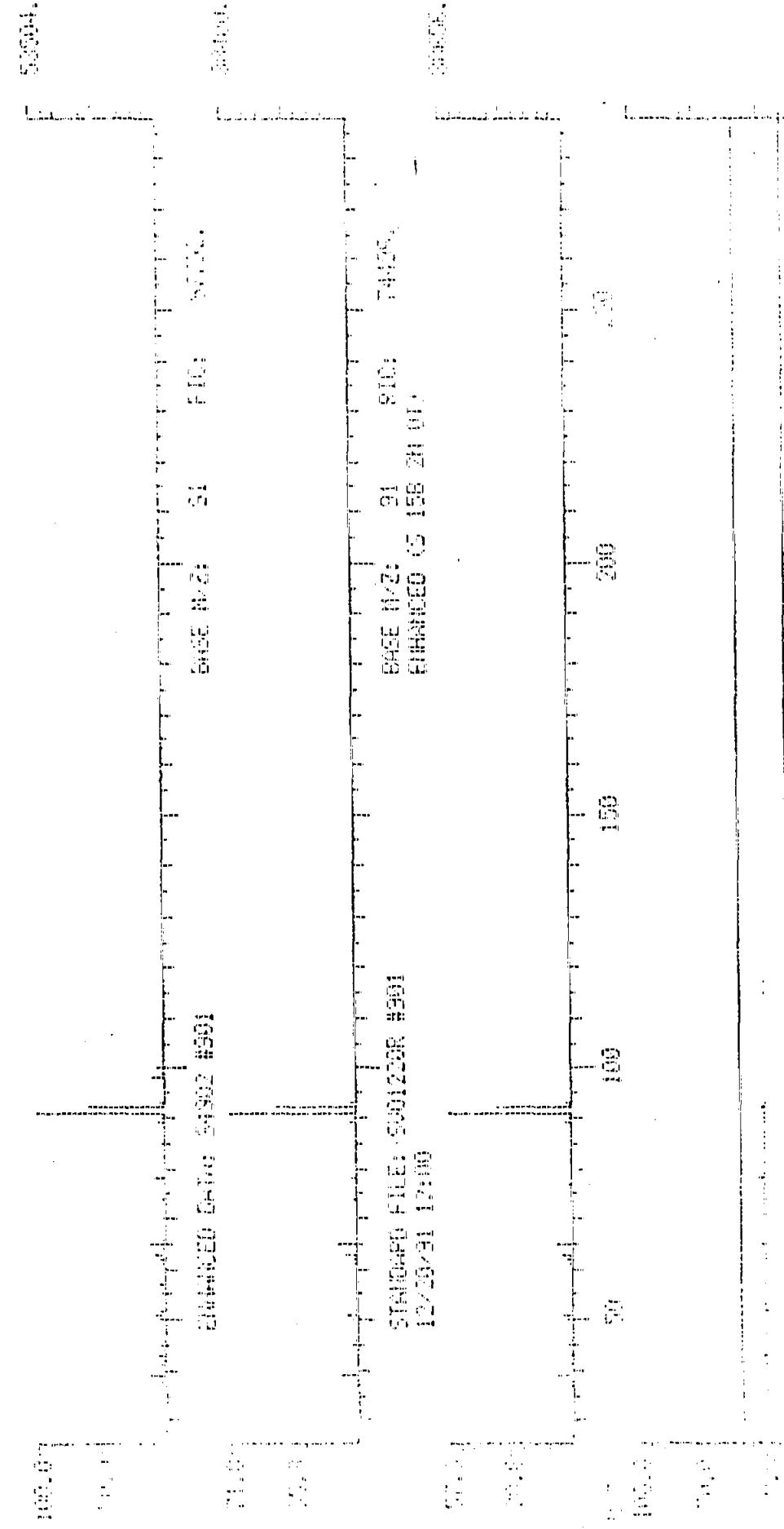
AR100709

00077

DATA FILE: 54902 #901
1625 ET Compound Dispersion
CONTINUO: 1230 TO LINE
12/26/91 16:38

STANDARD FILE: Standard #901
CONTINUO: 54902 #901

DATA NO: 91
PAGE NO: 91
PIC: 17000.



BR100710

00078

DATA FILE: 54962 #1014
TARGET COMPOUND COMPARISON
COMPOUND: C24H 5HEXENE

RAU DATA: 54962 #1014
12/20/91 16:36

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STANDARD FILE: SU01220R #1015
CALL: 54962 #3

RAU DATA: 54962 #1014
12/20/91 16:36

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600.0
400.0
200.0
0.0

1000.0
800.0
600.0
400.0
200.0
0.0

1000.0
800.0
600.0
400.0
200.0
0.0

MR100711

00079

DATA FILE: S4962 #1185
TARGET COMPOUND COMPARISON
COMPOUND: C150 0BP XYLENE

STANDARD FILE: S001220R #1185
CALL: S4962 #3

SAN FRANCISCO 54962 #1185
12/20/91 16:36
BASE M/Z: 91 RIC: 1015800.

1000.0
100.0
10.0
1.0
1.0E-01
1.0E-02
1.0E-03
1.0E-04
1.0E-05

ENHANCED 100.000 S4962 #1185
12/20/91 16:36
BASE M/Z: 91 RIC: 456728.

1000.0
100.0
10.0
1.0
1.0E-01
1.0E-02
1.0E-03
1.0E-04
1.0E-05

STANDARD FILE: S001220R #1185
12/20/91 16:36
BASE M/Z: 91 RIC: 500004.

1000.0
100.0
10.0
1.0
1.0E-01
1.0E-02
1.0E-03
1.0E-04
1.0E-05

STANDARD FILE: S001220R #1185
12/20/91 16:36
BASE M/Z: 91 RIC: 500004.
ENHANCED 0.5 15B 20.00

1000.0
100.0
10.0
1.0
1.0E-01
1.0E-02
1.0E-03
1.0E-04
1.0E-05

1000.0
100.0
10.0
1.0
1.0E-01
1.0E-02
1.0E-03
1.0E-04
1.0E-05

1000.0
100.0
10.0
1.0
1.0E-01
1.0E-02
1.0E-03
1.0E-04
1.0E-05

-50.0

-100.0

AR100712

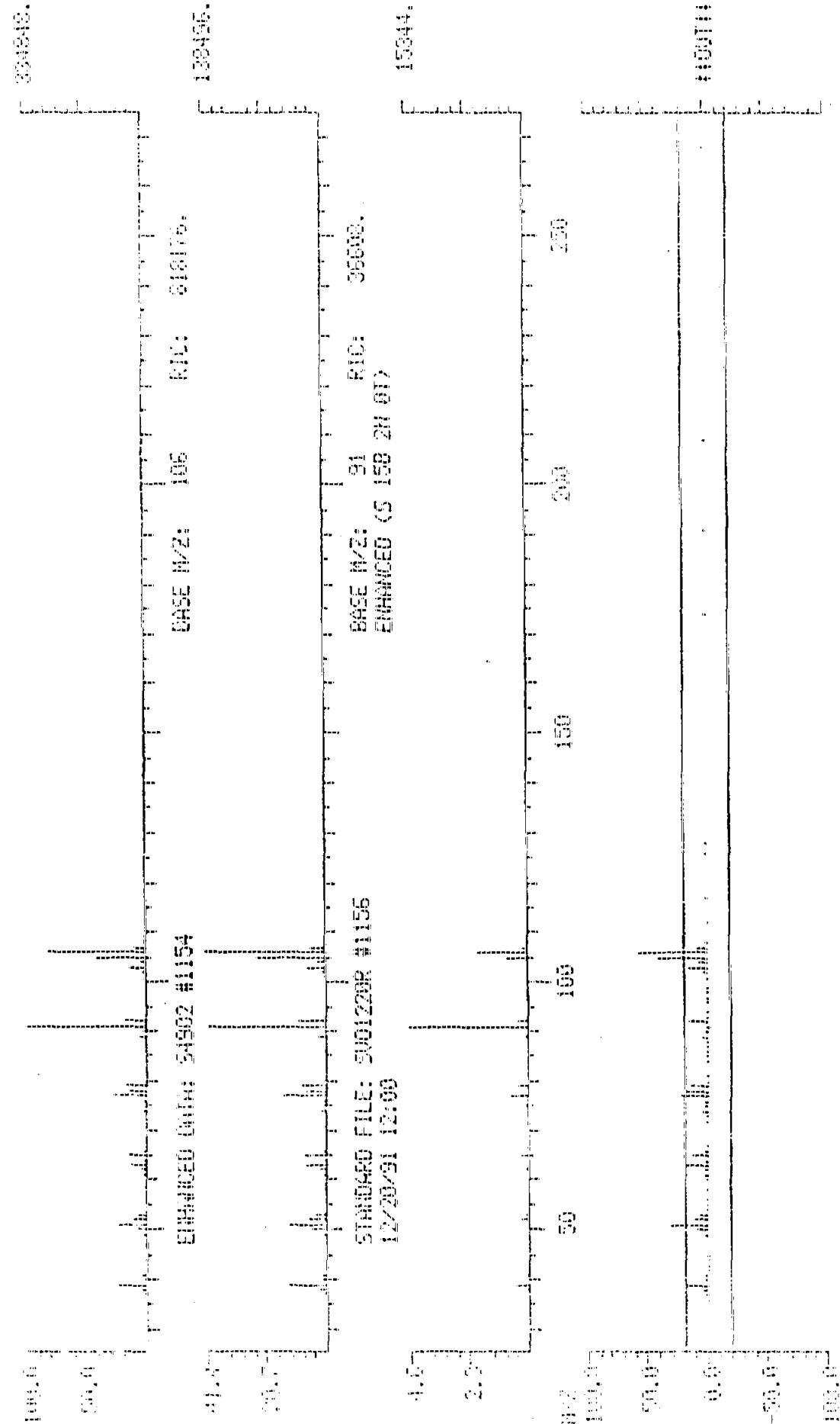
000.

DATA FILE: 54992 #1154
TARGET COMPARISON
COLLECT: C250 4-XYLENE

STANDARD FILE: 50012282 #1155
CALL: 54992 #15

RT# RT#C 54992 #1154
12/20/91 10:35

RT# RT#C 54992 #1155
12/20/91 10:35



20713

00000

CHROMATOGRAPHIC RECORDS FILE: 8400

Date: 8400, 71

Lab: 8400 190000-00

805 AM

Method: GLASS PLATE COLUMN METHOD 8400, C 10, 10 F:

Mobile: 100UL OF 100% V/VOL. Isopropanol HPLC

Submitted by: WESTON Analysts: CC

Revised by: G.G.C.
Date: 8400

REPRINT AREA: * USE ACTIVITY AREA & REG. PAGE
Repl. Fac.: from Library Entry

No.	Name
1	C101 BROMOCHLOROMETHANE ** INT. STD. **
2	C110 1,4-DIFLUORETENE ** INT. STD. **
3	C120 CHLOROBENZENE-DS ** INT. STD. **
4	C515 1,1,2-TRICHLOROETHANE-D4 ** S. STD. **
5	C800 TOLUENE-DS ** S. STD. **
6	C810 4-BROMOFLUORETENE ** S. STD. **
7	C910 CHLOROMETHANE **
8	C915 BROMINETHANE
9	C920 VINYL CHLORIDE *
10	C930 CHLOROETHANE
11	C940 METHYLENE CHLORIDE
12	C950 ACROLEIN
13	C960 ACETONE
14	C970 1,1,2,2-TETRA
15	C980 CARBON DISULFIDE
16	C985 1,1-DIFLUORETHENE *
17	C990 1,1-THIACHLOROETHANE **
18	C005 TRANS-1,2-DICHLOROETHENE
19	C000 TRICHLOROFLUOROMETHANE
20	C060 CHLOROFORM *
21	C065 1,1,2-TRICHLOROETHANE
22	C070 1,1,2,3-TETRA
23	C080 CHLOROACETIC ACID
24	C125 VINYL ACETATE
25	C130 BROMO-1,1-CHLOROMETHANE
26	C140 1,1,2-CHLOROPROPANE *
27	C145 TRANS-1,2-DICHLOROPROPENE
28	C150 1,1,2,3-TETRA
29	C160 1,1,2,3-TRICHLOROETHANE
30	C165 BENZENE
31	C170 1,1,2-TRICHLOROPROPENE
32	C175 2-CHLOROETHYL VINYL ETHER
33	C180 BROMOFORM **
34	C220 TETRACHLOROETHENE
35	C210 2-HEXANONE
36	C205 4-METHYL 2-PENTANONE
37	C225 1,1,2,3-TETRACHLOROETHANE ***
38	C230 TOLUENE *
39	C235 CHLOROBENZENE *
40	C440 ETHYL CHLORIDE *
41	C441 ETHYL CHLORIDE *
42	C442 ETHYL CHLORIDE *
43	C443 ETHYL CHLORIDE *

AR100114

00082

Sample
1000 CPE (AEM)

No.	RT/1	Scan	Time	Ref	RRT	Peak	Area(Height)	Amount	%Tot
1	49	337	8:26	1	1.000	A BB	114353.	50.000 NG	1.61
2	114	747	18:45	2	1.000	A BB	361966.	50.000 NG	1.61
3	117	937	23:30	3	1.000	A BB	285690.	50.000 NG	1.61
4	68	447	11:11	1	1.326	A BB	101240.	57.127 NG	2.07
5	98	591	22:18	3	0.949	A BB	320315.	52.481 NG	1.91
6	NOT FOUND	1094	Mortix Interference	cannot integrate					
7	70	51	11:17	1	0.151	A BB	2616	1.203 NG ✓	0.04
8	NOT FOUND								
9	NOT FOUND								
10	NOT FOUND								
11	49	227	8:38	1	0.668	A VB	4521.	5.282 NG ✓	0.12
12	NOT FOUND								
13	42	361	8:38	1	0.774	A BB	10234.	52.512 NG ✓	1.19
14	NOT FOUND								
15	NOT FOUND								
16	NOT FOUND								
17	NOT FOUND								
18	NOT FOUND								
19	NOT FOUND								
20	NOT FOUND								
21	NOT FOUND								
22	NOT FOUND								
23	NOT FOUND								
24	NOT FOUND								
25	95	627	18:42	2	0.837	A BB	46176	17.450 NG ✓	0.38
26	NOT FOUND								
27	NOT FOUND								
28	76	649	18:15	2	0.866	A BB	109364.	51.977 NG ✓	0.77
29	NOT FOUND								
30	NOT FOUND								
31	105	545	21:06	3	0.893	A BB	7703460.	294.000 100.000 ✓	25.85
32	NOT FOUND								
33	NOT FOUND								
34	91	824	22:26	3	0.957	A BB	23748.	2.350 NG FT	0.18
35	NOT FOUND								
36	105	1015	23:24	3	1.081	A BB	6510.	2.350 NG FT	0.18
37	NOT FOUND								
38	NOT FOUND								
39	NOT FOUND								
40	NOT FOUND								
41	105	1123	23:27	3	1.260	A BB	266680.	48.917 100.000 ✓	1.47

AR100715

00083

PC
12/20/91 19:02:00 ~~5.00~~
SAMPLE: 6360-2
CHROM.: GLASS PACKED COLUMN/METHOD 624/1.C 10/18/91
PULSE: 6 1,400 LABEL: N 0, 4.0 60MM A 0, 1.0 J 0 502: 0 26, 5
S42

1000

DATA: S4205 #48 SC405 39 TO 1400

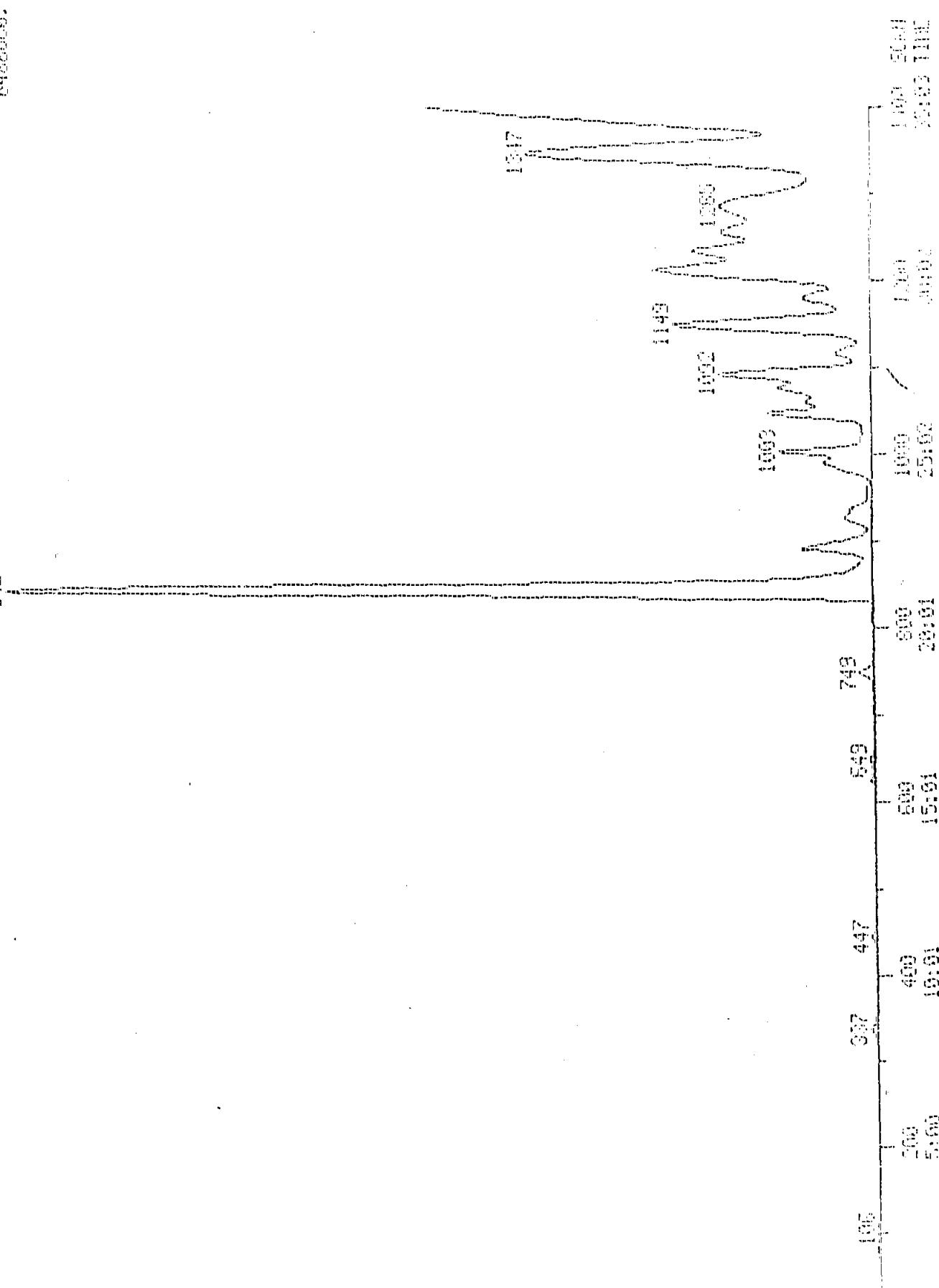
CALI: S4005 #3

SC405 39 TO 1400

S4205.

AR100716

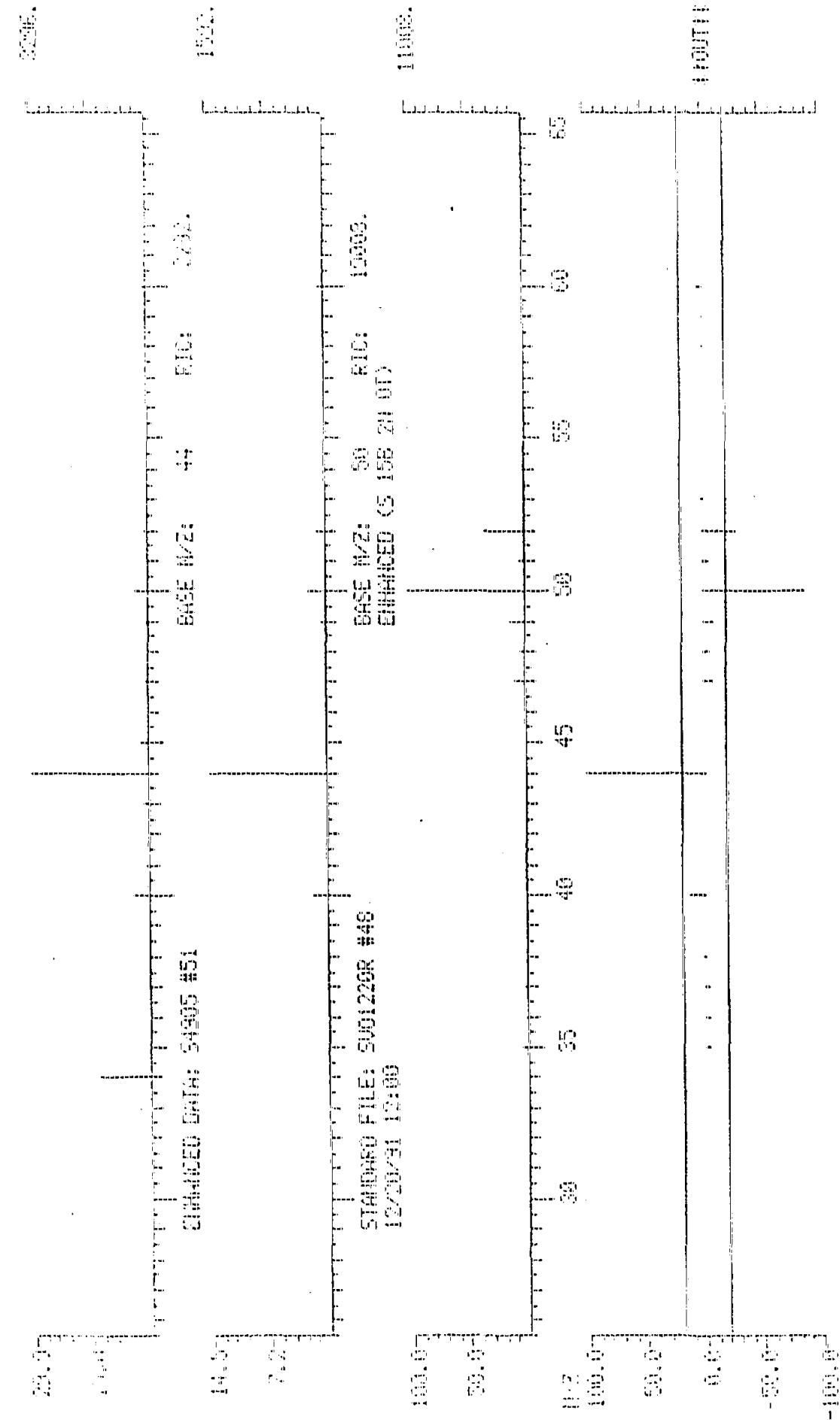
* 00084



DATA FILE: 54365 #51
TARGET COMPOUND COMPARISON
COMPOUND: 2010 CHLOROMETHANE
REF. DATA: 54365 #51
12/20/91 12:02

STANDARD FILE: S001226R #48
CALL: 54365 #8

REF. DATA: 54365 #51
12/20/91 12:02



AR100717

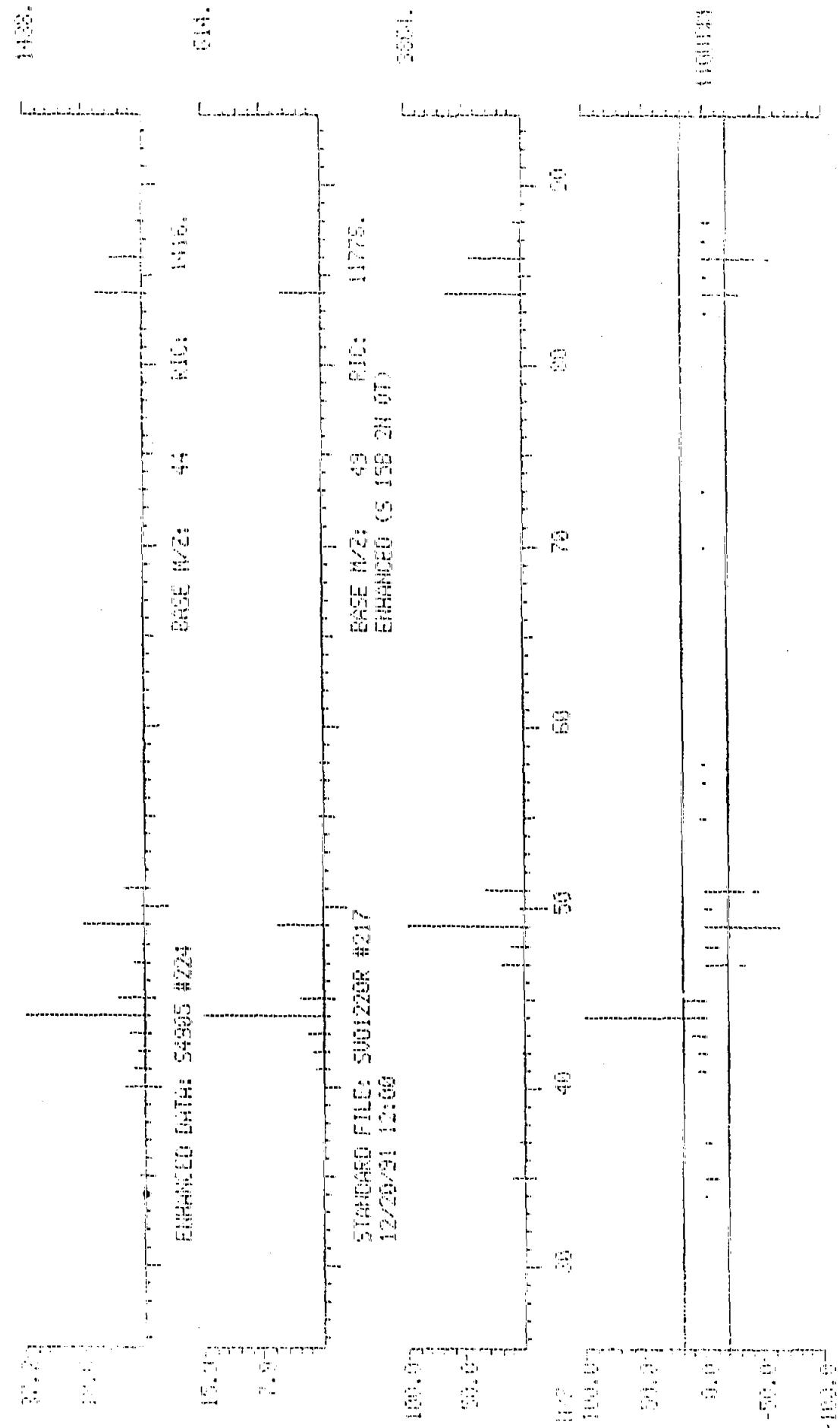
00085

DATA FILE: 54965 #224
TARGET COMPOUND COMPARISON
COMPOUND: 0639 METHYLENE CHLORIDE

STANDARD FILE: S001220R #217
COMPOUND: 54965 #3

REF DATE: 54965 #224
12/20/91 12:00

REF DATE: 54965 #224
12/20/91 12:00



AP-00718

00086

DATA FILE: S4995 #251
TARGET COMPOUND COMPARISON
COMPOUND: C635 ACETONE

STANDARD FILE: S001226R #259
CALI: S4995 #2

94% DATA: 54995 #251
12/26/91 19:30

89% DATA: 54995 #252

81% DATA: 54995 #253

74% DATA: 54995 #254

68% DATA: 54995 #255

62% DATA: 54995 #256

56% DATA: 54995 #257

50% DATA: 54995 #258

44% DATA: 54995 #259

38% DATA: 54995 #260

32% DATA: 54995 #261

26% DATA: 54995 #262

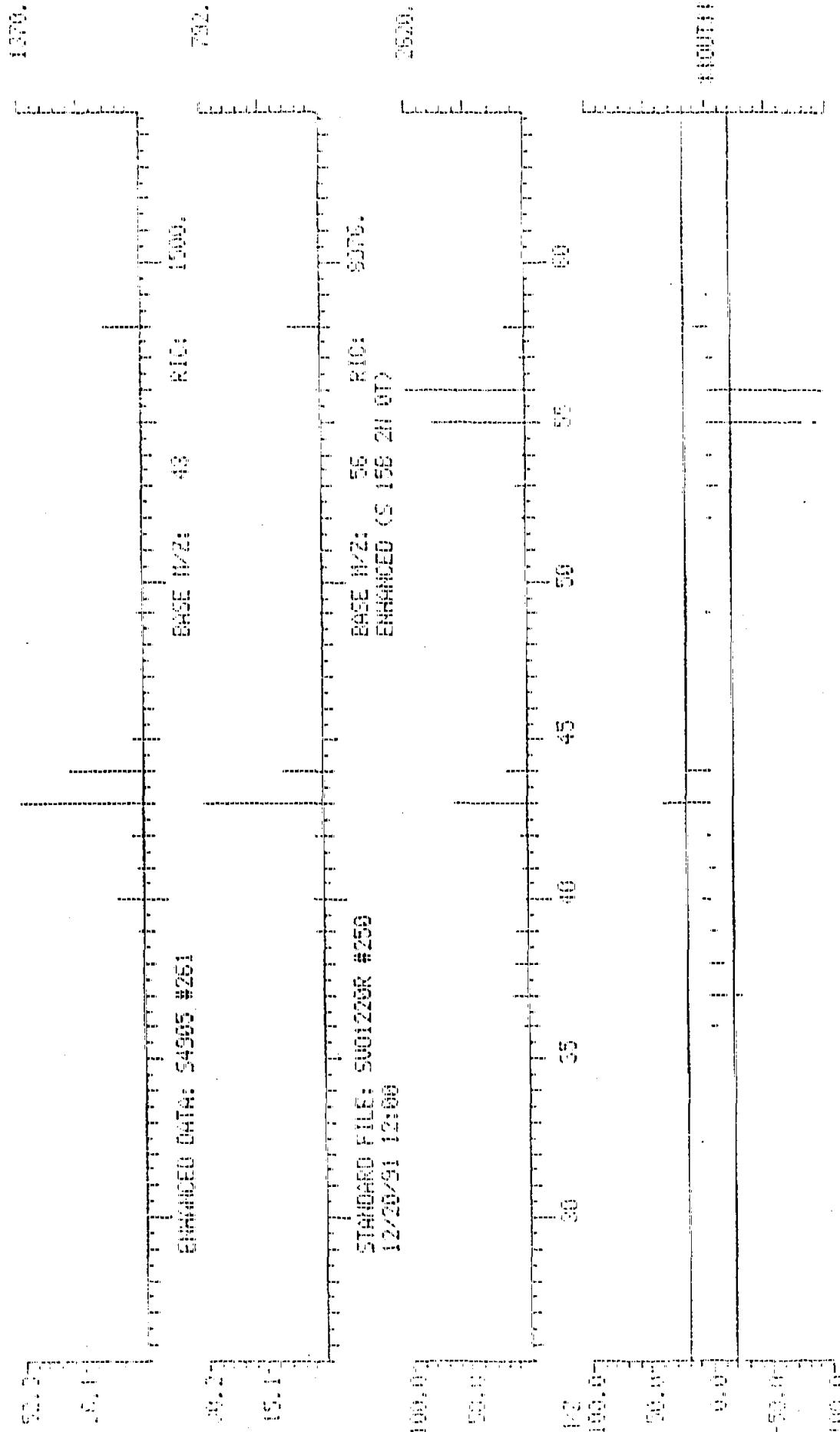
20% DATA: 54995 #263

14% DATA: 54995 #264

8% DATA: 54995 #265

2% DATA: 54995 #266

0% DATA: 54995 #267



PR 100719

00087

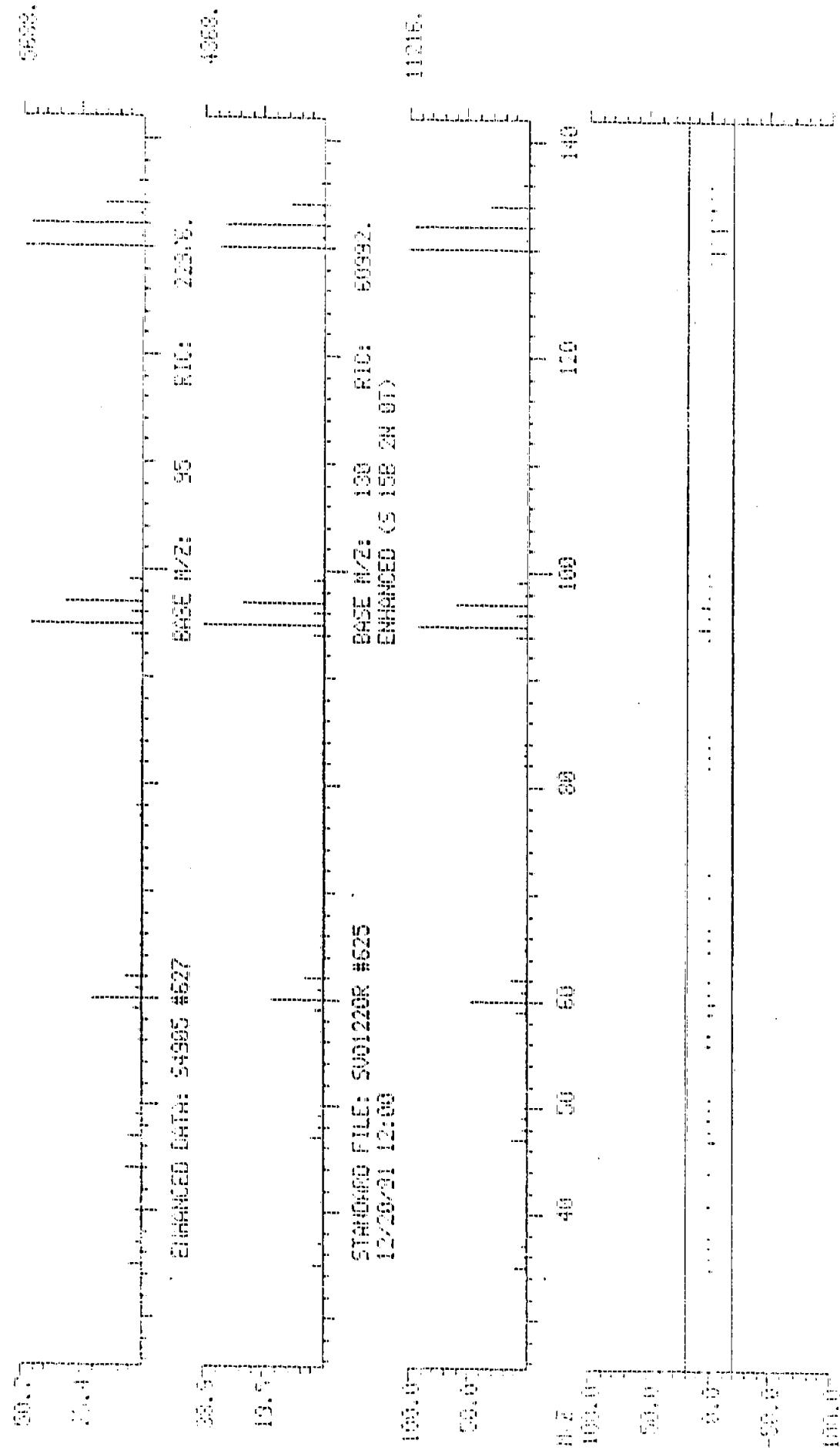
DATA FILE: S4905 #627

STANDARD FILE: S491220R #625

TARGET COMPOUND COMPARISON
COMPOUND: C130 TRICHLOROETHENE

RAW DATA: 54905 4527
12/26/91 12:02

BASE M/Z: 130 RIC: 31994.



AR100720

00083

DATA FILE: 54905 #649
TARGET COMPOUND COMPARISON
COMPOUND: C16S BENZENE

STANDARD FILE: SU01220R #648
CALL: 54905 #6

SPM DATE: 12/26/91 12:00

BASE M/Z: 78 SIC: 30172.

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ENHANCED SPECTRUM 54905 #649

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STANDARD FILE: SU01220R #648

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PR10072

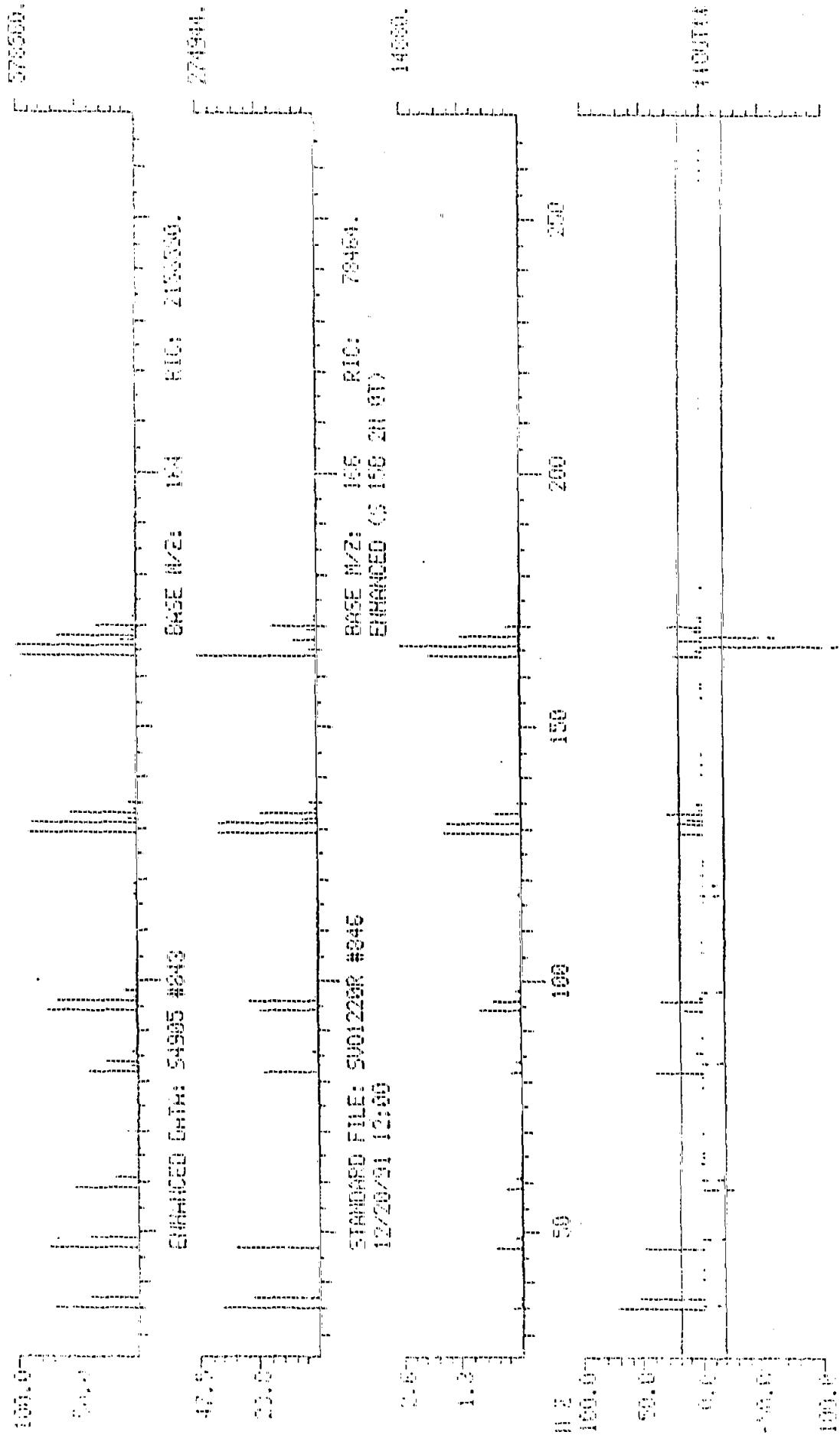
00089

DATA FILE: S4995 #845
TARGET COMPOUND COPPER(II)
COMPOUND: C229 TETRAOCHLOROTHEINE

STANDARD FILE: S00122008 #845
CALL #: S4995 #8

PROBE DATE: 12/20/01 19:02
SAMPLE #: S4995 #845

REF ID: S4995 #8



BR100722

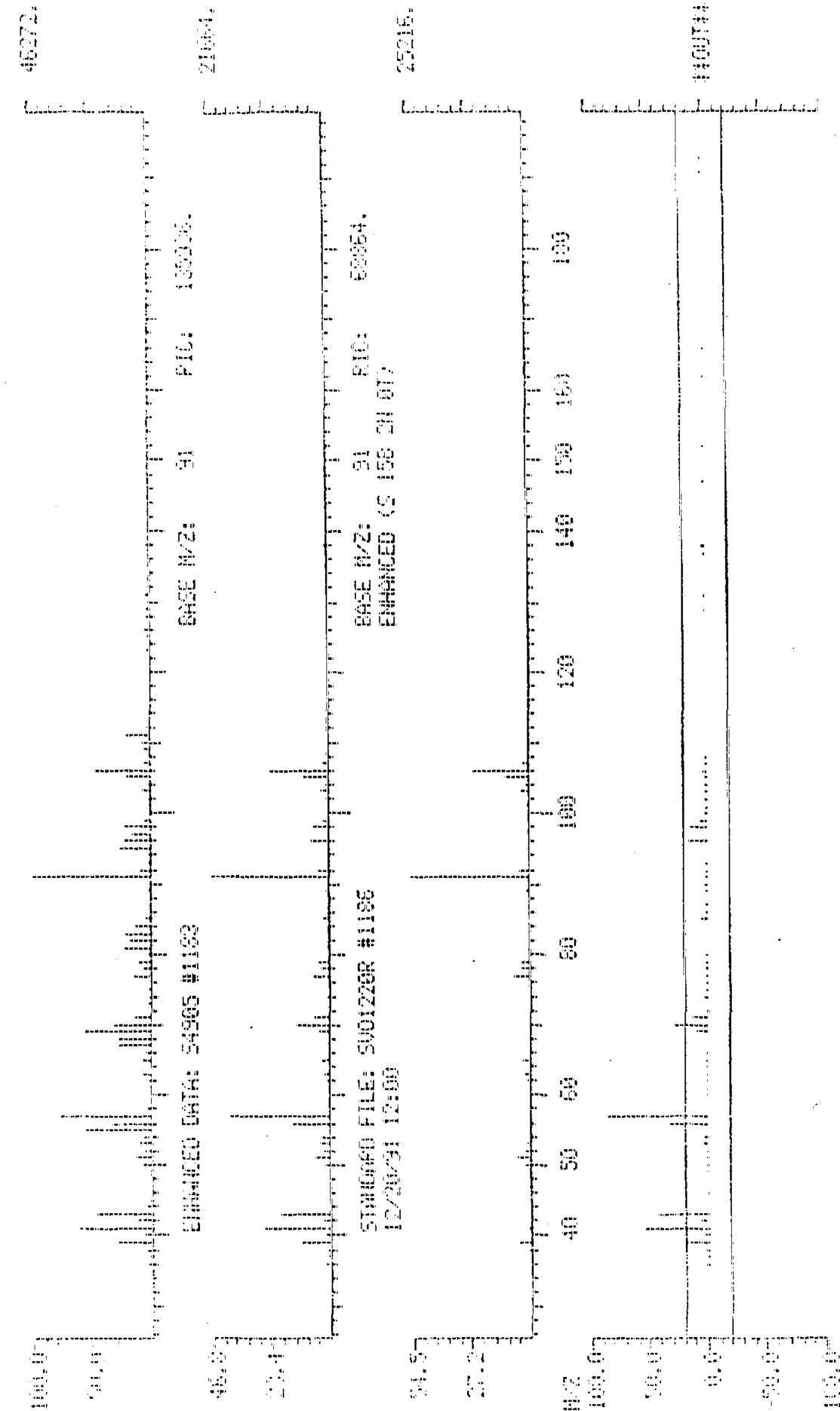
00090

DATA FILE: 54905 #1185
 TARGET COMPOUND: C256 OMP XYL ENE
 COMPOUND #: C256 OMP XYL ENE
 RUN DATE: 12/20/91 15:00Z
 S4905 #1185

STANDARD FILE: S001228R #1185

CALI: 54905 #2

SIGC: 44905.



ER100723

00091

PRINCETON TESTING LABORATORY

Weston TAT
5 Underwood Court
Delran, New Jersey 08075

Job Number: 9108360-002GMWH
Report Date: 12/16/91

BULK ASBESTOS ANALYSIS REPORT - METHOD: EPA-600/M4-82-020

SAMPLE IDENTIFICATION:	001 A07	
ANALYTICAL METHOD:	Polarized Light Microscopy With Dispersion Staining	
GROSS SAMPLE APPEARANCE:	Homogeneous Fibrous Gray	
SAMPLE TREATMENT:	Homogenized	
ASBESTOS FIBERS PRESENT?:	Yes	
ASBESTOS CONCENTRATION:		
AMOSITE	45%	
CHRYSOTILE	0%	
CROCIDOLITE	0%	
OTHER	0%	
TOTAL	45%	
OTHER FIBROUS MATERIAL:		
FIBROUS GLASS		
CELLULOSE		
OTHER		
NONFIBROUS MATERIAL PRESENT:	Silicates Binders Others	
ANALYST:	Susan Sharples	

AR100724

00092