



SDMS DocID 259202

Superfund records Center
SITE: Davis Liquid
BREAK: 3.2
OTHER: _____

VOLUME 4
PRIVATE WELL
LABORATORY RESULTS
For
Davis Liquid Site

REM II

**PERFORMANCE OF REMEDIAL RESPONSE
ACTIVITIES AT UNCONTROLLED
HAZARDOUS WASTE SITES**

U.S. EPA CONTRACT NO. 68-01-6939

**CAMP DRESSER & MCKEE INC.
ROY F. WESTON, INC.
WOODWARD-CLYDE CONSULTANTS
CLEMENT ASSOCIATES, INC.
ICF INCORPORATED
C. C. JOHNSON & ASSOCIATES, INC.**

VOLUME 4
PRIVATE WELL
LABORATORY RESULTS
For
Davis Liquid Site

EPA Contract No.: 58-01-6939/117/WP1

Work Assignment No.: 18-I1107

Document No.: 117-WP1-RT-BJMQ-1

DAVIS LIQUID SITE
Smithfield, Rhode Island

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SERIES: Private Wells

<u>Case Number</u>	<u>Volume</u>
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INTRODUCTION

This volume is apart of a series that presents all validated organic laboratory analyses results pertaining to the private well sampling conducted at the Davis Liquid Site in Smithfield, Rhode Island. The series is a collection of five different cases in which each case is contained in two volumes. The first case, Number 3633, was received by Radian Laboratory on December 7, 1984 and included 14 samples. The second case, Number 3969, was received by GCA Corporation on March 1, 1985 and included 15 samples. The third case, Number 3981, was received by GCA Corporation on March 8, 1985 and included 15 samples. The fourth case, Number 4005, was received by ChemTech on March 13, 1985 and included 10 private well samples. The final case, Number 4124, was received by GCA Corporation on April 5, 1985 and included 10 private well samples. This volume contains case Number 3969.

In addition to the lab results, two summary tables have been included. Table 1 is a list of each case number with the corresponding Sample Management Office (SMO) organic traffic numbers. Table 2 contains a summary of the validated lab analysis reports from the case contained in this volume. At the top of the summary sheets, the well location and resident are matched to their respective SMO traffic numbers.

Organic analyses results from each case are each contained in bound volumes marked accordingly on the covers. These volumes have been tabbed at the beginning of each sample by their sample SMO traffic numbers for easy reference. The first page of results, at the tab are the volatile compounds, the second page contains the semi-volatile compounds and the third page contains the pesticide/PCB compounds. These three pages contain the compounds on the Hazardous Substances List (HSL). Immediately following are two pages which allow write-in space for an additional 10 volatile compounds and 20 semi-volatile compounds which may be matched through a mass spectral library search. They are listed as tentatively present. The rest of the volume consists of the lab QA/QC program.

NOTE:

The validation process has not yet been completed for the fourth and fifth cases, numbers 4005 and 4124. Upon receiving validation, these two cases will be added to this collection of data making this series on private well data complete. Validation completion is expected by mid-September, 1985.

TABLE 1

PRIVATE WELLS

CASE #3969 AB 025
AB 026
AB 027
AB 028
AB 029
AB 030
AB 031
AB 032
AB 033
AB 034
AB 035
AB 036
AB 037
AB 038
AB 039

4

TABLE 2

3/8/85

Volatile Organics Analysis

PRIVATE WELLS

CAS Number	Sample ID EPA SMO	TIMENTA	HAYWARD	CRETELLA	DUP TO PW-03	McANN	BASTIEN	CASEY	DUP TO PW-07	TRIP BLANK	MOWRY	TIERNEY
		AB 025	AB 026	AB 027	AB 028	AB 029	AB 030	AB 031	AB 032	AB 033	AB 034	AB 035
		PW-07	PW-06	PW-05	PW-04	PW-03	PW-02	PW-01	PW-08	PW-15	PW-12	PW-13
74-87-3	Chloromethane											
74-83-9	Bromomethane											
75-01-4	Vinyl Chloride											
75-00-3	Chloroethane											
75-09-2	Methylene Chloride	R	R	R	R	R	R	R	R	R	R	R
67-64-1	Acetone	46										
75-15-0	Carbon Disulfide				R							
75-35-4	1, 1-Dichloroethene											
75-34-3	1, 1-Dichloroethane											
156-60-5	Trans-1, 2-Dichloroethene											
67-66-3	Chloroform				13 K							
107-06-2	1, 2-Dichloroethane											
78-93-3	2-Butanone											
71-55-6	1, 1, 1-Trichloroethane											
56-23-5	Carbon Tetrachloride											
108-05-4	Vinyl Acetate											
75-27-4	Bromodichloromethane											
79-34-5	1, 1, 2, 2-Tetrachloroethane											
78-87-5	1, 2-Dichloropropane											
10061-02-6	Trans-1, 3-Dichloropropene											
79-01-6	Trichloroethene											
124-48-1	Dibromochloromethane											
79-00-5	1, 1, 2-Trichloroethane											
71-43-2	Benzene					7.3						
10061-01-5	cis-1, 3-Dichloropropene											
110-75-8	2-Chloroethylvinylether											
75-25-2	Bromoform											
591-78-6	2-Hexanone											
108-10-1	4-Methyl-2-Pentanone											
127-18-4	Tetrachloroethene											
108-88-3	Toluene											
108-90-7	Chlorobenzene											
100-41-4	Ethylbenzene											
100-42-5	Styrene											
	Total Xylenes						14 J					

Total ppm's = 46 0 0 8.9 7.3 104 7 0 0 0 0

Semivolatile Organics Analysis

Sample ID		AB 025	AB 026	AB 027	AB 028	AB 029	AB 030	AB 031	AB 032	AB 033	AB 034	AB 35
CAS Number	EPA SMO											
62-75-9	N-Nitrosodimethylanine											
108-95-2	Phenol											
62-53-3	Aniline											
111-44-4	bis(2-Chloroethyl)Ether											
95-57-8	2-Chlorophenol											
541-73-1	1,3-Dichlorobenzene											
106-46-7	1,4-Dichlorobenzene											
100-51-6	Benzyl Alcohol											
95-50-1	1,2-Dichlorobenzene											
95-48-7	2-Methylphenol											
39638-32-9	bis(2-chloroisopropyl)Ether											
106-44-5	4-Methylphenol											
621-64-7	N-Nitroso-Di-n-Propylamine											
67-72-1	Hexachloroethane											
98-95-3	Nitrobenzene											
78-59-1	Isophorone											
88-75-5	2-Nitrophenol											
105-67-9	2,4-Dimethylphenol											
65-85-0	Benzoic Acid											
111-91-1	bis(2-Chloroethoxy)Methane	12	R	R	R	R	R	R	R	R	R	R
120-83-2	2,4-Dichlorophenol											
120-82-1	1,2,4-Trichlorobenzene											
91-20-3	Naphthalene											
106-47-8	4-Chloroaniline											
87-68-3	Hexachlorobutadiene											
59-50-7	4-Chloro-3-Methylphenol											
91-57-6	2-Methylnaphthalene											
77-47-4	Hexachlorocyclopentadiene											
88-08-2	2,4,6-Trichlorophenol											
95-95-4	2,4,5-Trichlorophenol											
91-58-7	2-Chloronaphthalene											
88-74-4	2-Nitroaniline											
131-11-3	Dimethyl Phthalate											
208-98-8	Acenaphthylene											
99-09-2	3-Nitroaniline											

CAS Number	Sample ID EPA SMO	AB 25	AB 26	AB 27	AB 28	AB 29	AB 30	AB 31	AB 32	AB 33	AB 34	AB 35
		83-32-9	Acenaphthene									
51-28-5	2,4-Dinitrophenol											
100-02-7	4-Nitrophenol											
132-64-9	Dibenzofuran											
121-14-2	2,4-Dinitrotoluene											
608-20-2	2,6-Dinitrotoluene											
84-66-2	Diethylphthalate											
7005-72-3	4-Chlorophenyl-phenylether											
86-73-7	Fluorene							FK				
100-01-6	4-Nitroaniline											
534-52-1	4,6-Dinitro-2-Methylphenol											
86-30-6	N-Nitrosodiphenylamine (1)											
101-55-3	4-Bromophenyl-phenylether											
118-74-1	Hexachlorobenzene											
87-86-5	Pentachlorophenol											
85-01-8	Phenanthrene											
120-12-7	Anthracene											
84-74-2	Di-n-Butylphthalate											
206-44-0	Fluoranthene											
92-87-5	Benzidine											
129-00-0	Pyrene											
85-68-7	Butylbenzylphthalate											
91-94-1	3,3'-Dichlorobenzidine											
56-55-3	Benzo[a]Anthracene											
117-81-7	bis(2-Ethylhexyl)Phthalate								R R	R R	R R	R R
218-01-9	Chrysene											
117-84-0	Di-n-Octyl Phthalate											
205-99-2	Benzo[b]Fluoranthene											
207-06-9	Benzo[k]Fluoranthene											
50-32-8	Benzo[a]Pyrene											
193-39-5	Indeno[1,2,3-cd]Pyrene											
53-70-3	Dibenzo[a,h]Anthracene											
191-24-2	Benzo[g,h,i]Perylene											
	XYLENE							TP				

W/E # 3981
3/2/85

Case # 398
3/1/85

Volatile Organics Analysis

HOME
Sample ID
EPA
SMO

FIELD BLANK

OXILVIE

CAS
Number

CAS Number	Sample ID	COLPACK	HORNKR	TRIP BLANK	DUP TO PW-27	POWERS	MOORE	NYHAN	QUIMETTE	OSW
		AB050	AB051	AB054	AB052	AB053	AB036	AB037	AB038	AB039
		PW26	PW27	PW28	PW29	PW30	PW12	PW11	PW10	PW09
74-87-3	Chloromethane									
74-83-9	Bromomethane									
75-01-4	Vinyl Chloride	R	R	R	R	10				
75-00-3	Chloroethane									
75-09-2	Methylene Chloride	R	R	R	R	R	R	R	R	R
67-64-1	Acetone	R	R	R	R	11	R			
75-15-0	Carbon Disulfide									
75-35-4	1, 1-Dichloroethane			5 J						
75-34-3	1, 1-Dichloroethane									
156-60-5	Trans-1, 2-Dichloroethane									
67-66-3	Chloroform									
107-06-2	1, 2-Dichloroethane									
78-93-3	2-Butanone									
71-55-6	1, 1, 1-Trichloroethane									
56-23-5	Carbon Tetrachloride									
108-05-4	Vinyl Acetate									
75-27-4	Bromodichloromethane									
79-34-5	1, 1, 2, 2-Tetrachloroethane									
78-87-5	1, 2-Dichloropropane									
10061-02-6	Trans-1, 3-Dichloropropene									
79-01-6	Trichloroethene									
124-48-1	Dibromochloromethane									
79-00-5	1, 1, 2-Trichloroethane									
71-43-2	Benzene									
10061-01-5	cis-1, 3-Dichloropropene									
110-75-8	2-Chloroethylvinylether									
75-25-2	Bromoform	R	R	R	R	R				
591-78-6	2-Hexanone									
108-10-1	4-Methyl-2-Pentanone									
127-18-4	Tetrachloroethene									
108-88-3	Toluene									
108-90-7	Chlorobenzene									
100-41-4	Ethylbenzene									
100-42-5	Styrene	R	R	R	R	R				
	Total Xylenes									

Total ppm =

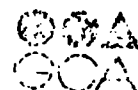
21 0 5 0 0 0 0 0 0

CAS Number	Sample ID EPA SMO	AB 050	AB 051	AB 052	AB 053	AB 36	AB 37	AB 38	AB 39
52-75-9	N-Nitrosodimethylaniline								
108-95-2	Phenol								
52-53-3	Aniline								
111-44-4	bis(2-Chloroethyl)Ether								
95-57-8	2-Chlorophenol								
541-73-1	1,3-Dichlorobenzene								
106-46-7	1,4-Dichlorobenzene								
100-51-6	Benzyl Alcohol								
35-50-1	1,2-Dichlorobenzene								
35-48-7	2-Methylphenol								
39638-32-9	bis(2-chloroisopropyl)Ether								
106-44-5	4-Methylphenol								
521-64-7	N-Nitroso-Di-n-Propylamine								
57-72-1	Hexachloroethane								
98-95-3	Nitrobenzene								
78-59-1	Isophorone								
88-75-5	2-Nitrophenol								
105-67-9	2,4-Dimethylphenol								
65-85-0	Benzoic Acid								
111-91-1	bis(2-Chloroethoxy)Methane								
120-83-2	2,4-Dichlorophenol								
120-82-1	1,2,4-Trichlorobenzene	R	R		R	R			
91-20-3	Naphthalene								
106-47-8	4-Chloroaniline								
87-68-3	Hexachlorobutadiene	R	R		R	R			
59-56-7	4-Chloro-3-Methylphenol								
91-57-6	2-Methylnaphthalene								
77-47-4	Hexachlorocyclopentadiene								
88-06-2	2,4,6-Trichlorophenol								
95-95-4	2,4,5-Trichlorophenol								
91-58-7	2-Chloronaphthalene								
88-74-4	2-Nitroaniline								
131-11-3	Dimethyl Phthalate								
208-96-8	Acenaphthylene	R	R		R	R			
99-09-2	3-Nitroaniline								

Semivolatile Organics Analysis

CAS Number	Sample ID EPA SMO	AB050	AB052	AB054	AB052	AB053	AG 30	AB 31	AB 38	AB 39
83-32-9	Acenaphthene									
51-28-5	2, 4-Dinitrophenol	R	R		R	R				
100-02-7	4-Nitrophenol									
132-64-9	Dibenzofuran									
121-14-2	2, 4-Dinitrotoluene									
606-20-2	2, 6-Dinitrotoluene									
84-66-2	Diethylphthalate									
7005-72-3	4-Chlorophenyl-phenylether									
86-73-7	Fluorene									
100-01-6	4-Nitroaniline									
534-52-1	4, 6-Dinitro-2-Methylphenol									
86-30-6	N-Nitrosodiphenylamine (1)									
101-55-3	4-Bromophenyl-phenylether									
118-74-1	Hexachlorobenzene									
87-86-5	Pentachlorophenol									
85-01-8	Phenanthrene									
120-12-7	Anthracene									
84-74-2	Di-n-Butylphthalate	R	R		R	R				
206-44-0	Fluoranthene									
92-87-5	Benzidine									
129-00-0	Pyrene									
85-68-7	Butylbenzylphthalate									
91-94-1	3, 3'-Dichlorobenzidine									
56-55-3	Benzo(a)Anthracene									
117-81-7	bis(2-Ethylhexyl)Phthalate	R	R		R	R	R	R	R	R
218-01-9	Chrysene									
117-84-0	Di-n-Octyl Phthalate	21								
205-99-2	Benzo(b)Fluoranthene									
207-08-9	Benzo(k)Fluoranthene									
50-32-8	Benzo(a)Pyrene									
193-39-5	Indeno(1, 2, 3-cd)Pyrene									
53-70-3	Dibenzo(a, h)Anthracene									
191-24-2	Benzo(g, h, i)Perylene									

IV. STANDARDS YACKET



GOA CORPORATION
Technology Division



Contractor GCA

Contract # 68-01-6767

CROSS REFERENCE TABLE - CASE 3969

1. Pesticides/PCBs

Date of Analysis of Standards

03/28/85 to 03/29/85

Applicable Samples

All samples primary column

AB 027, AB 028, AB 029, AB 032,
AB 034, AB 026 on primary
column

EXHIBIT C

Hazardous Substance List (HSL) and
Contract Required Detection Limits (CRDL)**

Volatiles	CAS Number	Detection Limits*	
		Low Water ^a ug/L	Low Soil/Sediment ^b ug/Kg
1. Chloromethane	74-87-3	10	10
2. Bromomethane	74-83-9	10	10
3. Vinyl Chloride	75-01-4	10	10
4. Chloroethane	75-00-3	10	10
5. Methylene Chloride	75-09-2	5	5
6. Acetone	67-64-1	10	10
7. Carbon Disulfide	75-15-0	5	5
8. 1,1-Dichloroethene	75-35-4	5	5
9. 1,1-Dichloroethane	75-35-3	5	5
10. trans-1,2-Dichloroethene	156-60-5	5	5
11. Chloroform	67-66-3	5	5
12. 1,2-Dichloroethane	107-06-2	5	5
13. 2-Butanone	78-93-3	10	10
14. 1,1,1-Trichloroethane	71-55-6	5	5
15. Carbon Tetrachloride	56-23-5	5	5
16. Vinyl Acetate	108-05-4	10	10
17. Bromodichloromethane	75-27-4	5	5
18. 1,1,2,2-Tetrachloroethane	79-34-5	5	5
19. 1,2-Dichloropropane	78-87-5	5	5
20. trans-1,3-Dichloropropene	10061-02-6	5	5
21. Trichloroethene	79-01-6	5	5
22. Dibromochloromethane	124-48-1	5	5
23. 1,1,2-Trichloroethane	79-00-5	5	5
24. Benzene	71-43-2	5	5
25. cis-1,3-Dichloropropene	10061-01-5	5	5

(continued)

IV

Volatiles	CAS. Number	Detection Limits*	
		Low Water ^a ug/L	Low Soil/Sediment ^b ug/Kg
26. 2-Chloroethyl Vinyl Ether	110-75-8	10	10
27. Bromoform	75-25-2	5	5
28. 2-Hexanone	591-78-6	10	10
29. 4-Methyl-2-pentanone	108-10-1	10	10
30. Tetrachloroethene	127-18-4	5	5
31. Toluene	108-88-3	5	5
32. Chlorobenzene	108-90-7	5	5
33. Ethyl Benzene	100-41-4	5	5
34. Styrene	100-42-5	5	5
35. Total Xylenes		5	5

^aMedium Water Contract Required Detection Limits (CRDL) for Volatile HSL Compounds are 100 times the individual Low Water CRDL.

^bMedium Soil/Sediment Contract Required Detection Limits (CRDL) for Volatile HSL Compounds are 100 times the individual Low Soil/Sediment CRDL.

C-2

IV 2

9/84 Rev

CONTRACT DETECTION LIMITS - SEMIVOLATILE COMPOUNDS

Laboratory Name: GCA Case No.: 2969 Concentration LOW

Sample Matrix: water Contract No.: 68-01-6767

CAS #	Compound	ug/l
87-68-3	hexachlorobutadiene	10
77-47-4	hexachlorocyclopentadiene	10
78-59-1	isophorone	10
91-20-3	naphthalene	10
98-95-3	nitrobenzene	10
62-75-9	N-nitrosodimethylamine	10
86-30-6	N-nitrosodiphenylamine	10
621-64-7	N-nitrosodipropylamine	10
117-81-7	bis(2-ethylhexyl)phthalate	10
85-68-7	benzyl butyl phthalate	10
84-74-2	di-n-butyl phthalate	10
117-84-0	di-n-octyl phthalate	10
84-66-2	diethyl phthalate	10
131-11-3	dimethyl phthalate	10
56-55-3	benzo(a)anthracene	10
50-32-8	benzo(a)pyrene	10
205-99-2	benzo(b)fluoranthene	10
207-08-9	benzo(k)fluoranthene	10
218-01-9	chrysene	10
208-96-8	acenaphthylene	10
120-12-7	anthracene	10
191-24-2	benzo(ghi)perylene	10
86-73-7	fluorene	10
85-01-8	phenanthrene	10
53-70-3	dibenzo(a,h)anthracene	10
193-39-5	indeno(1,2,3-cd)pyrene	10
129-00-0	pyrene	10
62-53-3	aniline	10
100-51-6	benzyl alcohol	10
106-47-8	4-chloroaniline	10
132-64-9	dibenzofuran	10
91-57-6	2-methylnaphthalene	10
88-74-4	2-nitroaniline	50
99-09-2	3-nitroaniline	50
100-01-6	4-nitroaniline	50

IV 3



GCA CORPORATION
Technology Division

22

CONTRACT DETECTION LIMITS -- SEMIVOLATILE COMPOUNDS

Laboratory Name: GCA Case No.: 3969 Concentration low

Sample Matrix: water Contract No.: 68-01-6767

CAS #	Compound	ug/
88-06-2	2,4,6-trichlorophenol	10
59-50-7	p-chloro-m-cresol	10
95-57-8	2-chlorophenol	10
120-83-2	2,4-dichlorophenol	10
105-67-9	2,4-dimethylphenol	10
88-75-5	2-nitrophenol	10
100-02-7	4-nitrophenol	50
51-28-5	2,4-dinitrophenol	50
534-52-1	4,6-dinitro-2-methylphenol	50
87-86-5	pentachlorophenol	50
108-95-2	phenol	10
65-85-0	benzoic acid	50
95-48-7	2-methylphenol	10
108-39-4	4-methylphenol	10
95-95-4	2,4,5-trichlorophenol	50
83-32-9	acnaphthene	10
92-87-5	benzidine	50
120-92-1	1,2,4-trichlorobenzene	10
118-74-1	hexachlorobenzene	10
67-72-1	hexachloroethane	10
111-44-4	bis(2-chloroethyl)ether	10
91-58-7	2-chloronaphthalene	10
95-50-1	1,2-dichlorobenzene	10
341-73-1	1,3-dichlorobenzene	10
106-46-7	1,4-dichlorobenzene	10
91-94-1	3,3'-dichlorobenzidine	20
121-14-2	2,4-dinitrotoluene	10
606-20-2	2,6-dinitrotoluene	10
122-66-7	1,2-diphenylhydrazine	10
206-44-0	fluoranthene	10
7005-72-3	4-chlorophenyl phenyl ether	10
101-55-3	4-bromophenyl phenyl ether	10
39638-32-9	bis(2-chloroisopropyl)ether	10
111-91-1	bis(2-chloroethoxy)methane	10

IV 4



GCA CORPORATION
Technology Division

CONTRACT DETECTION LIMITS - PESTICIDES

Pesticides	CAS Number	Detection Limits*	
		Low Water ^e ug/L	Low Soil/Sediment ^f ug/Kg
104. alpha-BHC	319-84-6	0.05	8.0
105. beta-BHC	319-85-7	0.05	8.0
106. delta-BHC	319-86-8	0.05	8.0
107. gamma-BHC (Lindane)	58-89-9	0.05	8.0
108. Heptachlor	76-44-8	0.05	8.0
109. Aldrin	309-00-2	0.05	8.0
110. Heptachlor Epoxide	1024-57-3	0.05	8.0
111. Endosulfan I	959-98-8	0.05	8.0
112. Dieldrin	60-57-1	0.10	16.0
113. 4,4'-DDE	72-55-9	0.10	16.0
114. Endrin	72-20-8	0.10	16.0
115. Endosulfan II	33213-65-9	0.10	16.0
116. 4,4'-DDD	72-54-8	0.10	16.0
117. Endrin Aldehyde	7421-93-4	0.10	16.0
118. Endosulfan Sulfate	1031-07-8	0.10	16.0
119. 4,4'-DDT	50-29-3	0.10	16.0
120. Endrin Ketone	53494-70-5	0.10	16.0
121. Methoxychlor	72-43-5	0.5	80.0
122. Chlordane	57-74-9	0.5	80.0
123. Toxaphene	8001-35-2	1.0	160.0
124. AROCLOR-1016	12674-11-2	0.5	80.0
125. AROCLOR-1221	11104-28-2	0.5	80.0
126. AROCLOR-1232	11141-16-5	0.5	80.0
127. AROCLOR-1242	53469-21-9	0.5	80.0
128. AROCLOR-1248	12672-29-6	0.5	80.0
129. AROCLOR-1254	11097-69-1	1.0	160.0
130. AROCLOR-1260	11096-82-5	1.0	160.0

^eMedium Water Contract Required Detection Limits (CRDL) for Pesticide HSL Compounds are 100 times the individual Low Water CRDL.

^fMedium Soil/Sediment Contract Required Detection Limits (CRDL) for Pesticide HSL compounds are 15 times the individual Low Soil/Sediment CRDL.

*Detection limits listed for soil/sediment are based on wet weight. The detection limits calculated by the laboratory for soil/sediment, calculated on dry weight basis, as required by the contract, will be higher.

** Specific detection limits are highly matrix dependent. The detection limits listed herein are provided for guidance and may not always be achievable.

IV 5

INITIAL CALIBRATION DATA
VOLATILE HSL COMPOUNDS
(PAGE 1)

CASE NO. 3931,3969
CONTRACTOR GCA
CONTRACT NO. 68-01-6767

INSTRUMENT ID FINNIGAN
CALIBRATION DATE 3/3/85

MINIMUM RF FOR SPCC IS .300 MAXIMUM % RSD FOR CCC IS 30%

LABORATORY ID

COMPOUND	RF20	RF50	RF100	RF150	RF200	RF	%RSD	CCC* SPCC**
CHLOROMETHANE	1	1.14	1.1	1.23	1.32	1.15	8	**
BROMOMETHANE	1.2	1.27	1.21	1.17	1.2	1.21	1	
VINYL CHLORIDE	1.2	1.16	1.07	1.04	1.03	1.09	5	*
CHLOROETHANE	.86	.92	.84	.79	.81	.84	4	
METHYLENE CHLORIDE	1.8	1.87	1.6	1.6	1.53	1.67	7	
ACETONE	.19	.14	.23	.27	.23	.21	17	
CARBON DISULPHIDE	3.1	3.68	3.3	3.23	3.31	3.32	4	
1,1-DICHLOROETHANE	1.21	1.26	1.2	1.2	1.21	1.21	1	*
1,1-DICHLOROETHANE	2.9	3.37	3	3.3	3.4	3.19	6	**
TRANS-1,2-DICHLOROETHENE	1.6	1.65	1.5	1.5	1.5	1.55	3	
CHLOROFORM	4.3	4.46	4	4.3	4.46	4.3	2	*
1,2-DICHLOROETHANE	3.6	4.1	4	4.23	4.3	4.04	4	
2-BUTANONE	^	1.16	1.26	1.4	.84	1.16	14	
1,1,1-TRICHLOROETHANE	3.5	4	3.7	3.9	4.1	3.84	5	
CARBON TETRACHLORIDE	2.4	2.7	2.4	2.7	2.9	2.62	6	
VINYL ACETATE	.04	.06	.21	.21	.13	.13	49	
BROMODICHLOROMETHANE	.89	.92	.93	1.02	.97	.94	4	
1,2-DICHLOROPROPANE	.4	.41	.41	.46	.43	.42	4	*
TRANS-1,3-DICHLOROPROPENE	.36	.45	.47	.55	.55	.47	12	
TRICHLOROETHENE	.35	.35	.4	.43	.35	.37	8	
DIBROMOCHLOROMETHANE	.53	.54	.56	.62	.56	.56	4	
1,1,2-TRICHLOROETHANE	.27	.27	.31	.33	.28	.29	7	
BENZENE	1.1	1.17	1.2	1.2	1.2	1.17	2	
CIS-1,3-DICHLOROPROPENE	.18	.29	.34	.43	.45	.33	24	
2-CHLOROETHYL VINYLETHER	^	.01	.02	.01	.01	.01	30	
BROMOFORM	.34	.23	.38	.26	.21	.28	21	**
2-HEXANONE	^	.01	.19	.22	.1	.13	57	
4-METHYL-2-PENTANONE	^	.76	.89	.88	.6	.78	13	
TETRACHLOROETHENE	.5	.5	.46	.5	.47	.48	3	
1,1,2,2-TETRACHLOROETHANE	.3	.34	.4	.43	.45	.38	13	**
TOLUENE	1	1.02	1	.88	1.04	.98	4	*
CHLOROBENZENE	.91	.97	.93	1.1	.98	.97	5	**
ETHYLBENZENE	.44	.5	.46	.5	.49	.47	4	*
STYRENE	^	.17	.25	.34	.35	.27	24	
TOTAL XYLENES	1	1	1	1	1	1	0	

RESPONSE FACTOR (amount in ug/L) SPCC-SYSTEM PERFORMANCE CHECK COMPOUNDS (**)

RF-AVERAGE RESPONSE FACTOR

%RSD-PERCENT RELATIVE STANDARD DEVIATION

CCC-CALIBRATION CHECK COMPOUND (*)

10-5

Initial Calibration Data
Semivolatile HSL Compounds
 (Page 1)

Case No: 3969
 Contractor: GCM
 Contract No: 68-01-6767

Instrument ID: HP5985A
 Calibration Date: 3/25/85

Minimum RF for SPCC is 0.050 Maximum % RSD for CCC is 30%

Laboratory ID								
Compound	RF ₂₀	RF ₅₀	RF ₈₀	RF ₁₂₀	RF ₁₆₀	RF	% RSD	CCC- SPCC**
N-Nitrosodimethylamine								
Phenol	.70	.72	.84	.73	.74	0.75	7	*
Aniline	.83	.50	.71	.80	.72	0.81	8	
bis(2-Chloroethyl)Ether	.70	.60	.71	.60	.51	0.62	12	
2-Chlorophenol	.55	.57	.61	.60	.60	0.57	4	
1,3-Dichlorobenzene	.60	.53	.59	.52	.50	0.55	7	
1,4-Dichlorobenzene	.61	.52	.59	.55	.53	0.56	7	*
Benzyl Alcohol	.37	.34	.40	.34	.34	0.36	7	
1,2-Dichlorobenzene	.56	.52	.54	.49	.50	0.53	7	
2-Methylphenol	.54	.52	.67	.55	.62	0.58	4	
bis(2-chloroisopropyl)Ether	.52	.50	.53	.45	.45	0.47	7	
4-Methylphenol	.55	.60	.70	.60	.65	0.63	8	
N-Nitroso-Di-n-Propylamine	.50	.50	.44	.46	.44	0.45	5	**
Hexachloroethane	.20	.25	.24	.25	.25	0.27	8	
Nitrobenzene	.74	.70	.72	.71	.67	0.71	3	
Isophorone	1.23	1.25	1.52	1.27	1.20	1.31	8	
2-Nitrophenol	1.57	1.53	1.80	1.67	1.40	1.64	8	*
2,4-Dimethylphenol	2.24	2.51	2.60	2.71	2.60	2.55	5	
Benzoic Acid	†	.90	1.20	1.37	1.50	1.26	16	
bis(2-Chloroethoxy)Methane	1.00	1.61	1.81	1.55	1.70	1.54	8	
2,4-Dichlorophenol	1.75	1.80	2.05	2.15	2.00	1.97	17	
1,2,4-Trichlorobenzene	2.20	2.20	2.64	2.60	2.00	2.37	10	
Naphthalene	4.14	1.10	1.23	1.09	1.01	1.13	9	
4-Chloroaniline	3.70	3.43	4.60	4.20	3.90	4.07	8	
Hexachlorobutadiene	.90	1.00	1.14	1.06	1.00	1.07	7	*
4-Chloro-3-Methylphenol	1.20	1.60	2.20	2.04	2.00	1.87	17	*
2-Methylnaphthalene	2.25	2.20	2.60	2.30	2.00	2.24	9	
Hexachlorocyclopentadiene	.17	.16	.21	.18	.16	0.18	11	**
2,4,6-Trichlorophenol		.19	.24	.21	.21	0.21	8	*
2,4,5-Trichlorophenol	†	.19	.24	.21	.21	0.21	8	
2-Chloronaphthalene	.23	.21	.30	.24	.22	0.24	10	
2-Nitroaniline	†	.29	.40	.34	.33	0.34	12	
Dimethyl Phthalate	.08	.07	.08	.08	.07	0.08	6	
Acenaphthylene	.37	.24	.27	.24	.21	0.25	9	
3-Nitroaniline	†	.33	.29	.28	.30	0.33	10	
Acenaphthene	.41	.30	.44	.40	.36	0.39	8	*
2,4-Dinitrophenol	†	.07	.11	.11	.13	0.11	21	**
4-Nitrophenol	†	.20	.23	.20	.21	0.21	6	**
Dibenzofuran	.47	.41	.44	.42	.40	0.44	8	

Response Factor (subscript is the amount of nanograms)
 RF - Average Response Factor
 %RSD - Percent Relative Standard Deviation
 CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)
 † - not detectable at 20 ng

IV 7
 Form VI

**Initial Calibration Data
Semivolatile HSL Compounds
(Page 2)**

Case No: 3969
 Contractor: CCA
 Contract No: 68-01-6767

Instrument ID: HP5985A
 Calibration Date: 3/25/85

Minimum \bar{RF} for SPCC is 0.050 Maximum % RSD for CCC is 30%

Laboratory ID								
Compound	RF ₂₀	RF ₅₀	RF ₈₀	RF ₁₂₀	RF ₁₆₀	\bar{RF}	% RSD	CCC-SPCC**
2, 4-Dinitrotoluene	.30	.25	.32	.30	.20	0.29	8	
2, 6-Dinitrotoluene	.21	.20	.23	.22	.21	0.22	5	
Diethylphthalate	.16	.15	.17	.15	.14	0.15	7	
4-Chlorophenyl-phenylether	.33	.30	.36	.33	.31	0.33	6	
Fluorene	.80	.70	.82	.71	.70	0.75	7	
4-Nitroaniline	†	.25	.29	.29	.28	0.28	6	
4, 6-Dinitro-2-Methylphenol	†	.29	.13	.12	.11	0.11	13	
N-Nitrosodiphenylamine (1)	.13	.12	.14	.12	.12	0.13	6	*
4-Bromophenyl-phenylether	.10	.10	.11	.10	.10	0.10	4	
Hexachlorobenzene	.10	.10	.11	.10	.10	0.10	4	*
Pentachlorophenol	†	.65	.67	.67	.66	0.66	13	
Phenanthrene	.65	.64	.66	.66	.66	0.65	15	
Anthracene	.65	.64	.66	.66	.66	0.65	15	
Di-N-Butylphthalate	.10	.07	.10	.08	.09	0.09	14	
Fluoranthene	.66	.59	.70	.52	.50	0.57	13	*
Benzidine	†							**
Pyrene	2.00	1.70	2.26	2.02	2.50	2.16	17	
Butylbenzylphthalate	.92	.96	1.14	1.04	1.26	1.05	13	
3, 3'-Dichlorobenzidine								
Benzo(a)Anthracene	.30	.36	.43	.41	.41	0.38	12	
bis(2-Ethylhexyl)Phthalate	1.00	1.00	1.13	1.00	1.12	1.05	6	
Chrysene	.70	.36	.43	.40	1.71	0.35	12	
Di-n-Octyl Phthalate	3.51	4.40	5.97	3.50	3.95	4.01	27	*
Benzo(b)Fluoranthene	.91	1.05	1.31	1.90	0.86	1.01	16	
Benzo(k)Fluoranthene	.91	1.05	1.31	0.90	0.86	1.01	16	
Benzo(a)Pyrene	.71	.60	.74	0.65	0.62	0.66	8	*
Indeno(1, 2, 3-cd)Pyrene	.50	.26	.50	0.45	0.65	0.41	33	
Dibenz(a, h)Anthracene	.34	.15	.35	0.33	0.40	0.32	27	
Benzo(g, h, i)Perylene	.26	.20	.30	0.35	0.45	0.34	22	

Response Factor (subscript is the amount of nanograms)
 \bar{RF} - Average Response Factor
 %RSD - Percent Relative Standard Deviation
 CCC - Calibration Check Compounds (*)

SPCC - System Performance Check Compounds (**)
 † - not detectable at 20 ng
 (1) - Cannot be separated from diphenylamine

Form VI

IV 8

Contract No. 5-10-630 PM Column SP 2250/2401

Date of Analysis 3/29/85 Instrument ID HP 5870

EVALUATION CHECK FOR LINEARITY

LABORATORY ID	Eval. A	Eval. B	Eval. C	
PESTICIDE	CALIBRATION FACTOR EVAL. MIX A	CALIBRATION FACTOR EVAL. MIX B	CALIBRATION FACTOR EVAL. MIX C	% RSD (≤10%)
ALDRIN	7451/.025 277040	17613/.05 353760	46713/.125 375744	40152/54265 11.7 *
ENDRIN	10520/.05 210400	27427/.125 235432	62634/.25 248136	14201/231322 8.3
4,4'-DDT	5724/.05 116470	15772/.125 126176	4777/.25 191108	40570/144508 28.1 *
DIBUTYL CHLORENDATE	13443/.05 278860	34494/.125 275952	69680/.25 278720	16406/23474 6.57

EVALUATION CHECK FOR 4,4'-DDT/ENDRIN BREAKDOWN

	PERCENT BREAKDOWN EXPRESSED AS TOTAL DEGRADATION			
	EVAL. MIX B	EVAL. MIX B	EVAL. MIX B	EVAL. MIX B
ENDRIN	0	0	0	
4,4'-DDT	0	0	0	
LABORATORY ID	Eval. B	Eval. B	Eval. B	
TIME OF ANALYSIS	10:57 3/29	10:02 3/30	3:30 3/31	

EVALUATION OF RETENTION TIME SHIFT FOR DIBUTYLCHLORENDATE

SMO SAMPLE NO.	LAB ID	TIME OF ANALYSIS	PERCENT DIFF. *	SMO SAMPLE NO.	LAB ID	TIME OF ANALYSIS	PERCENT DIFF. *
Eval. A	Eval. A	10:00 3/29	31.74 RT				
Eval. B	Eval. B	10:57	0.03				
Eval. C	Eval. C	11:30	0.03				
Calibration	P-31-4	13:48	0.16				
AB 030	43220	5:26 3/30	0.27	BL			
Calibration	P-31-3	4:31 3/30	0.51				
AB 030	43220	5:26 3/30	0.27				
Eval. B	Eval. B	10:02	0.22				
Calibration	P-31-4	11:53	0.16				
Calibration	P-31-3	21:04	0.25				
Calibration	P-31-5	22:54	0.31				
Calibration	P-31-2	0:45 3/31	0.35				
Calibration	P-31-1	2:35	0.41				
Eval. B	Eval. B	3:30	0.44				

* ≤ 2% PACKED, 50.3% CAPILLARY.

4184

* See narrative

FORM VIII

IV 9

Contract No. 3-763-050 PV Column SP 2250/2401
 Date of Analysis 4/2/85 Instrument ID HP 5780

EVALUATION CHECK FOR LINEARITY

LABORATORY ID	Eval. A	Eval. B	Eval. C	
PESTICIDE	CALIBRATION FACTOR EVAL. MIX A	CALIBRATION FACTOR EVAL. MIX B	CALIBRATION FACTOR EVAL. MIX C	% RSD (≤10%)
ALDRIN	7537/.025 301480	17058/.05 341160	48527/.125 370616	34674/357752 10
ENDRIN	10273/.05 205460	25526/.125 204208	56511/.25 226364	12446/212010 5.9
4,4'-DDT	9727/.05 * 194540	21636/.125 173088	52098/.25 208392	17788/192007 9.3
DIBUTYL CHLORENDATE	15857/.05 317160	41436/.125, 125 331488 BL	77165/.25 308660	11537/319102 3.6

EVALUATION CHECK FOR 4,4'-DDT/ENDRIN BREAKDOWN

	PERCENT BREAKDOWN EXPRESSED AS TOTAL DEGRADATION			
	EVAL. MIX B	EVAL. MIX B	EVAL. MIX B	EVAL. MIX B
ENDRIN	0	0	0	0
4,4'-DDT	0	0	0	0
LABORATORY ID	Eval. B	Eval B	Eval B	Eval B
TIME OF ANALYSIS	12:46 4/1	3:34 4/2	9:53	16:59

EVALUATION OF RETENTION TIME SHIFT FOR DIBUTYLCHLORENDATE

SMO SAMPLE NO.	LAB ID	TIME OF ANALYSIS	PERCENT DIFF. *	SMO SAMPLE NO.	LAB ID	TIME OF ANALYSIS	PERCENT DIFF. *
Eval A	Eval A	11:51 4/1	31.67 RT				
Eval B	Eval B	12:46	0				
Eval C	Eval C	13:41	0.06				
Calibration	P-31-4	15:05	0.03				
ABO30 MS	CC-961MS	23:22	0.09				
Eval B	Eval B	3:34 4/2	0.06				
Calibration	P-31-4	4:59	0.12				
ABO34	43223	7:47	0				
Eval B	Eval B	9:53	0				
Calibration	P-31-5	12:05	0.03				
Calibration	P-31-3	13:29	0.03				
Calibration	P-31-2	14:53	0.09				
Calibration	P-31-1	16:17	0.32				
Eval B	Eval B	16:59	0.19				

* ≤ 2% PACKED, 30.3% CAPILLARY

4184

* See Narrative

FORM VIII

IV 10

Case No: CM 393 3969
 Contractor: LCC
 Contract No: 18-01-0077
 Instrument ID: Fujitsu Mat C-11A

Calibration Date: 3/4/85
 Time: 7:45
 Laboratory ID: V 2711
 Initial Calibration Date: 3/3/85

Minimum RF for SPCC is 0.300

Maximum %D for CCC is 25%

Compound	RF	RF50	%D	CCC	SPCC
Chloromethane	1.15	1.16	0.9		**
Bromomethane	1.21	1.28	5.1		
Vinyl Chloride	1.07	1.02	4.4	*	
Chloroethane	0.84	0.86	2.4		
Methylene Chloride	1.07	1.00	7.0		
Acetone	0.71	0.72	1.5		
Carbon Disulfide	2.22	2.00	11.0		
1,1-Dichloroethene	1.21	1.30	7.4	*	
1,1-Dichloroethane	2.19	2.60	13		**
Trans-1,2-Dichloroethene	1.95	1.78	15		
Chloroform	4.23	4.30	3.0	*	
1,2-Dichloroethane	1.14	2.43	15		
2-Butanone	1.15	1.25	7.4		
1,1,1-Trichloroethane	2.94	2.22	16		
Carbon Tetrachloride	2.42	1.91	27		
Vinyl Acetate	0.13	0.27	100		
Bromodichloromethane	0.74	0.67	24		
1,2-Dichloropropane	0.43	0.48	14	*	
Trans-1,3-Dichloropropene	0.47	0.43	8.5		
Trichloroethene	0.27	0.24	3.1		
Dibromochloromethane	0.56	0.61	31		
1,1,2-Trichloroethane	0.24	0.25	14		
Benzene	1.17	1.2	2.5		
cis-1,3-Dichloropropene	0.33	0.27	17		
2-Chloroethylvinylether	0.61	1.01	0		
Bromoform	2.25	2.15	1.4		**
2-Hexanone	0.13	0.17	33		
4-Methyl-2-Pentanone	0.75	0.58	74		
Tetrachloroethene	0.48	0.40	17		
1,1,2,2-Tetrachloroethane	0.32	0.27	22		**
Toluene	0.92	1.0	3.0	*	
Chlorobenzene	2.97	1.0	3.0		**
Ethylbenzene	0.47	0.51	3.5	*	
Styrene	0.27	0.25	7.8		
Total Xylenes	0.0	0.0	*		

RF₅₀ - Response Factor from daily standard file at 50 ug/l
 RF - Average Response Factor from initial calibration Form VI

%D - Percent Difference
 CCC - Calibration Check Compounds (*)
 SPCC - System Performance Check Compounds (**)

Form VII *ND - See Narrative

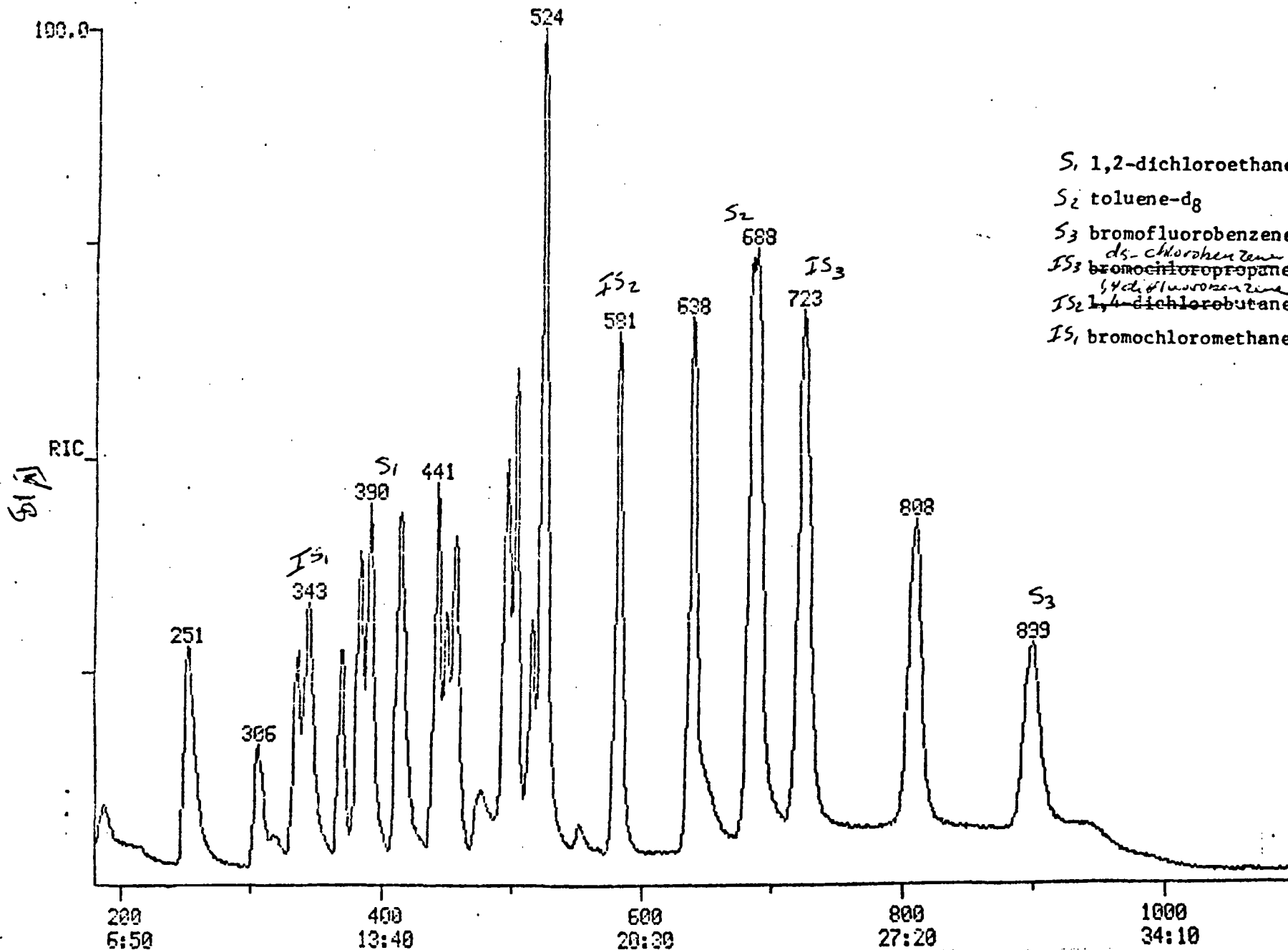
Handwritten signature/initials

RIC
03/04/85 7:45:00
SAMPLE: 50FF8 CASE 3931

DATA: U2716

SCANS 189 TO 1109

210176.



ORGANIC CHEMICALS IN WATER ANALYZER
CALIBRATION REPORT FILE: V2170

Case 3969
3451

DATE: V2170.TE
09/09 10:20:00
FILE:
SUBMITTED BY:

ANALYST:

UNIT-AREA(AIGHT) * REC. AREA / (REF. AREA(AIGHT) * RESP. FACT) *
P. FAC. FROM LIBRARY ENTRY

NO	NAME
1	BROMOCHLOROMETHANE (INTERNAL STANDARD 1)
2	CHLOROETHANE
3	BROMOMETHANE
4	VINYL CHLORIDE
5	CHLOROETHANE
6	METHYLENE CHLORIDE
7	ACETONE
8	CARBON DISULFIDE
9	1, 1-DICHLOROETHENE
10	1, 1-DICHLOROETHANE
11	TRANS-1, 2-DICHLOROETHENE
12	CHLOROFORM
13	CIS-1, 2-DICHLOROETHANE (SURROGATE 1)
14	1, 2-DICHLOROETHENE
15	2-BUTANONE
16	1, 1, 1-TRICHLOROETHANE
17	CARBON TETRACHLORIDE
18	VINYL ACETATE
19	BROMODICHLOROMETHANE
20	1, 2-DICHLOROPROPANE
21	TRANS-1, 2-DICHLOROPROPANE
22	TRICHLOROETHENE
23	DIBROMOCHLOROMETHANE
24	1, 1, 2-TRICHLOROETHANE
25	BENZENE
26	CIS-1, 3-DICHLOROPROPENE
27	2-CHLOROETHYL VINYL ETHER
28	1, 4-DIFLUOROBENZENE (INTERNAL STANDARD 2)
29	BROMOFORM
30	4-ETHYL-2-PENTANONE
31	1, 1, 2, 2-TETRACHLOROETHANE
32	TETRACHLOROETHENE
33	2-HEXANONE
34	DO-TOLUENE (SURROGATE 2)
35	TOLUENE
36	DS-CHLOROBENZENE (INTERNAL STANDARD 3)
37	CHLOROBENZENE
38	ETHYL BENZENE
39	4-BROMOFLUOROBENZENE (SURROGATE 3)
40	STYRENE
41	TOTAL XYLENES

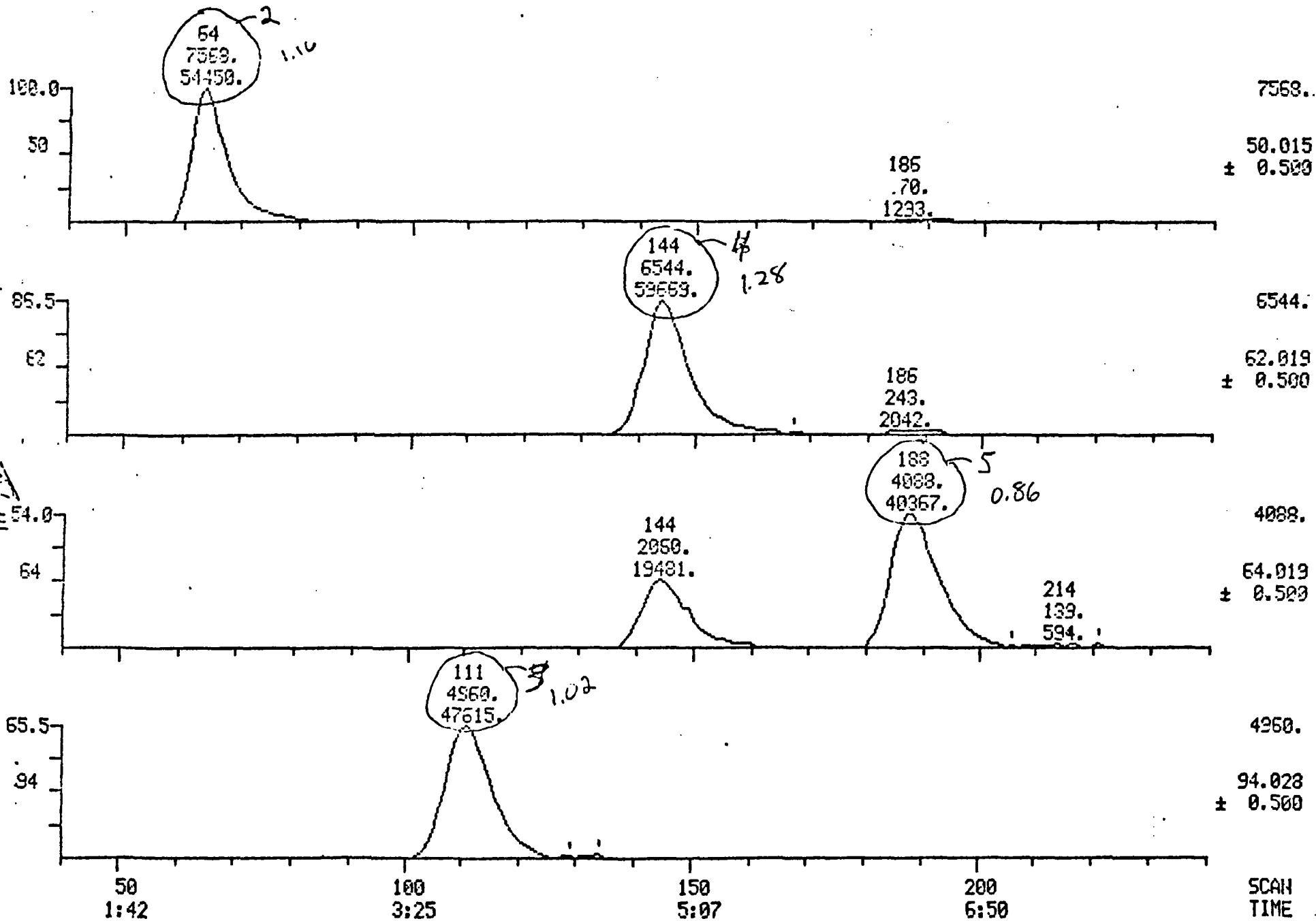
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IV 14

MASS CHROMATOGRAMS
03/04/85 7:45:00
SAMPLE: 50PPB CASE 3931

DATA: V2715

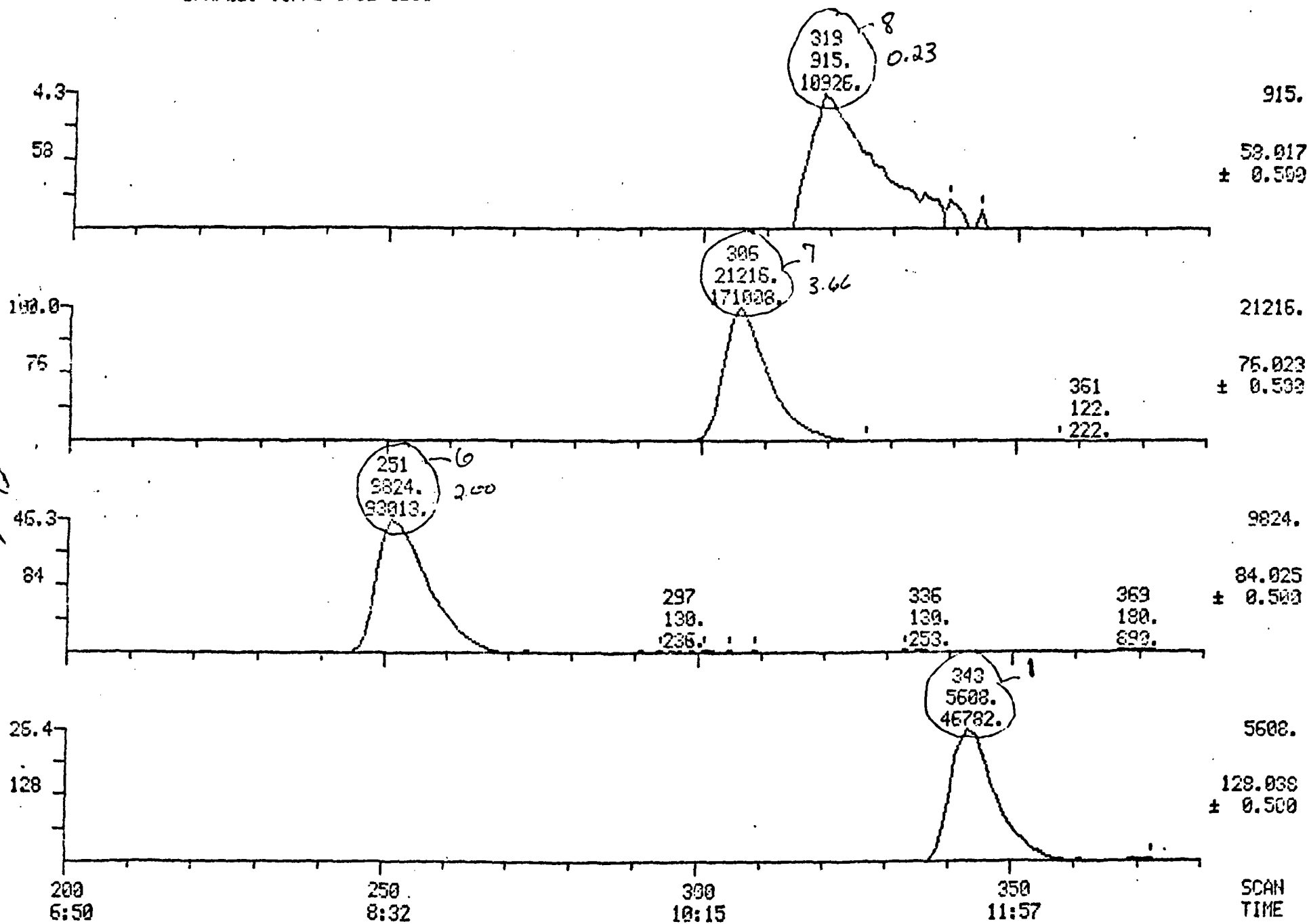
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03/04/85 7:45:00
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DATA: U2715

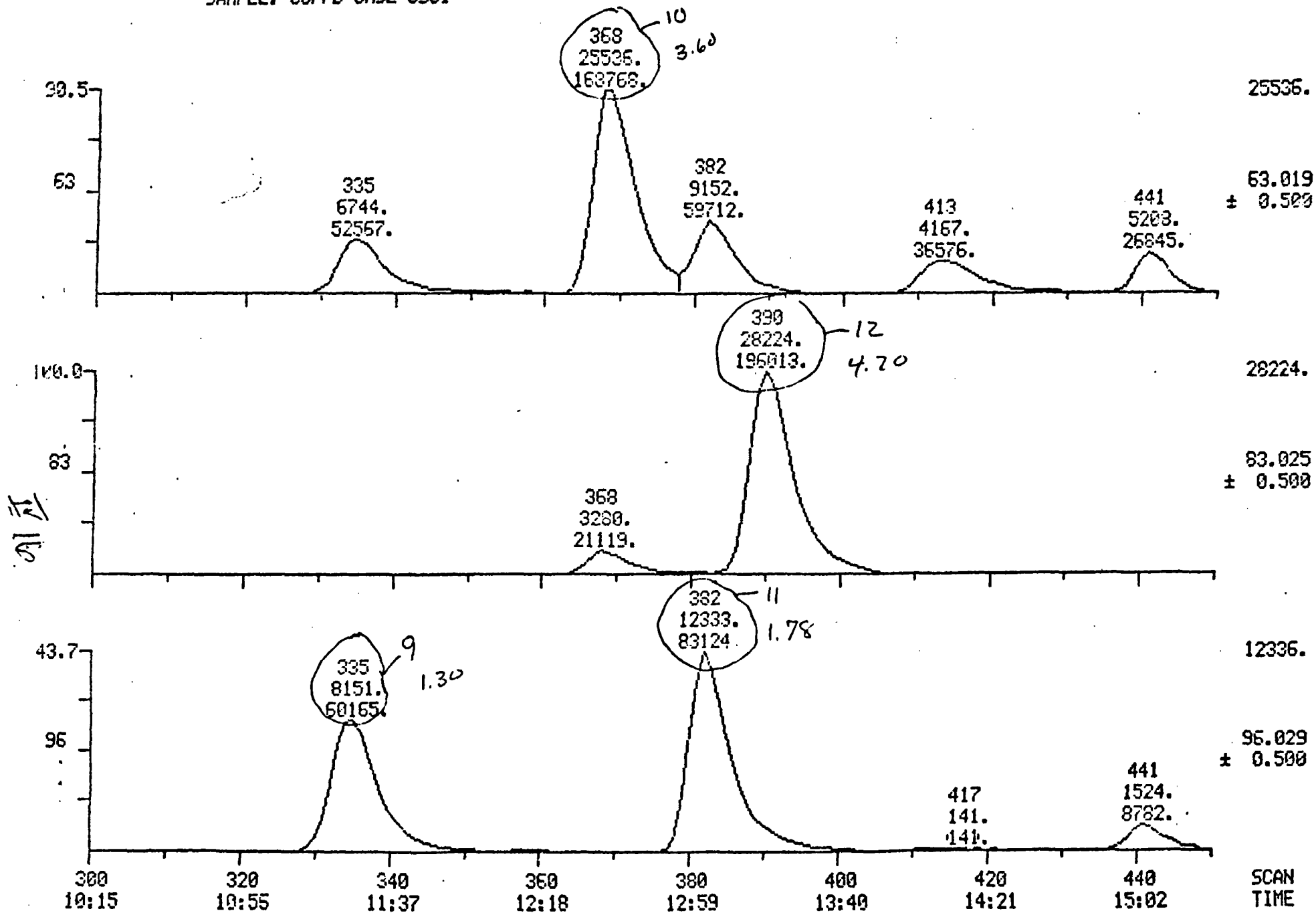
SCANS 200 TO 380



MASS CHROMATOGRAMS
 03/04/85 7:45:00
 SAMPLE: 50FPB CASE 3931

DATA: U2716

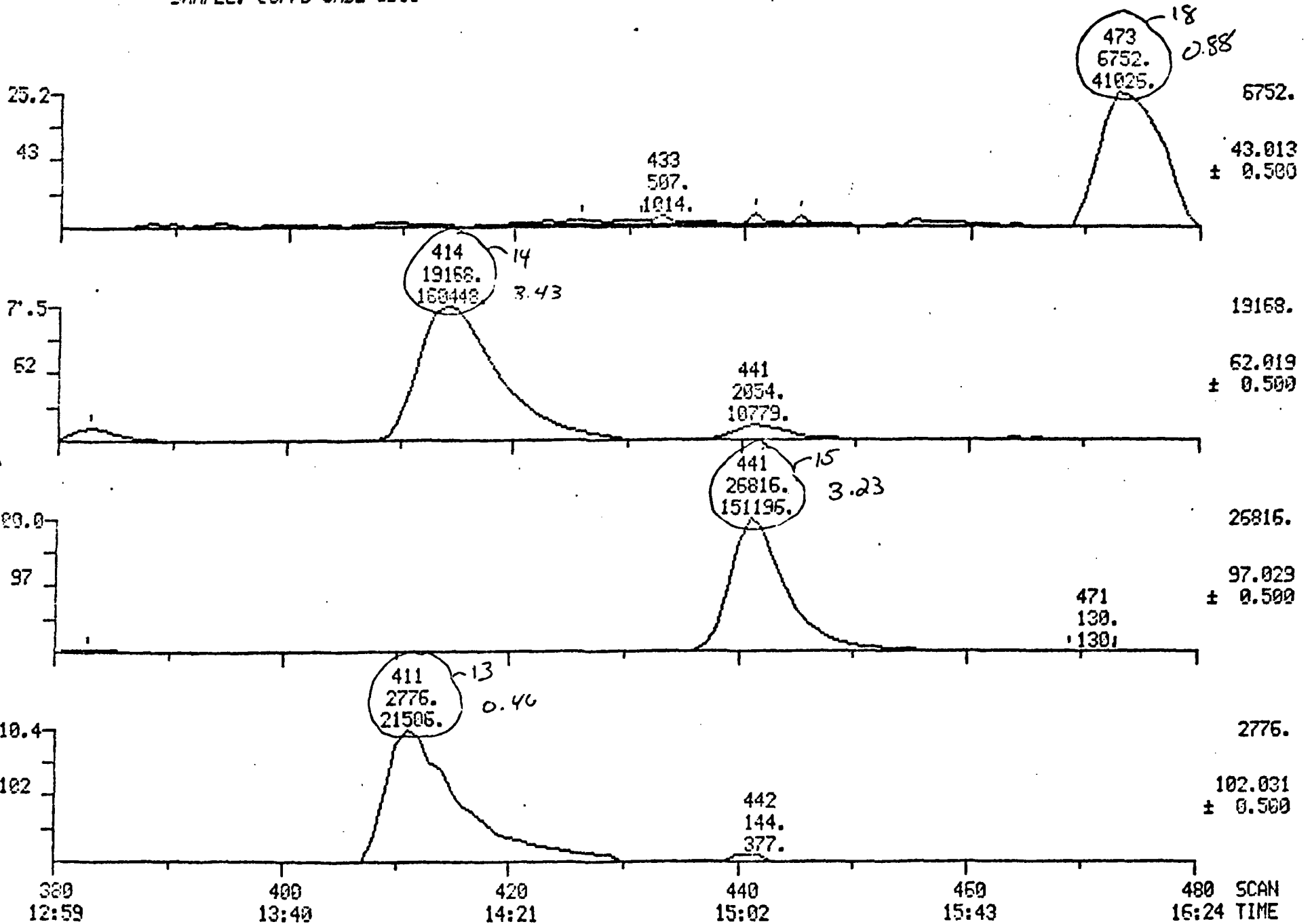
SCANS 300 TO 450



MASS CHROMATOGRAMS
03/04/85 7:45:00
SAMPLE: 50FPB CASE 3931

DATA: U2716

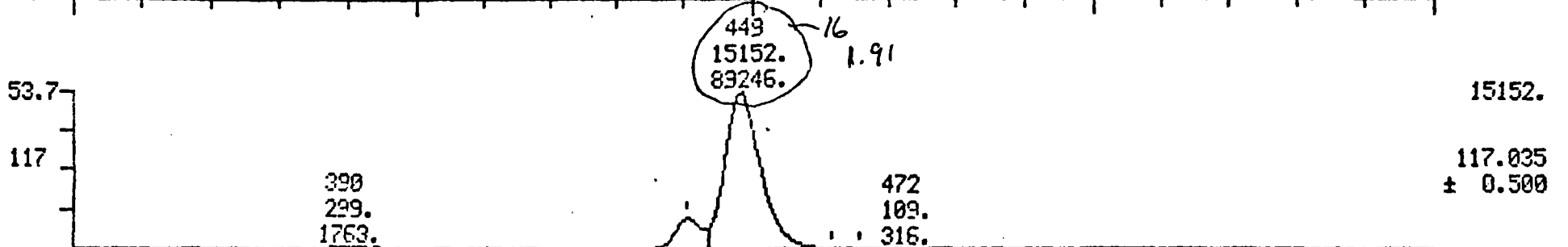
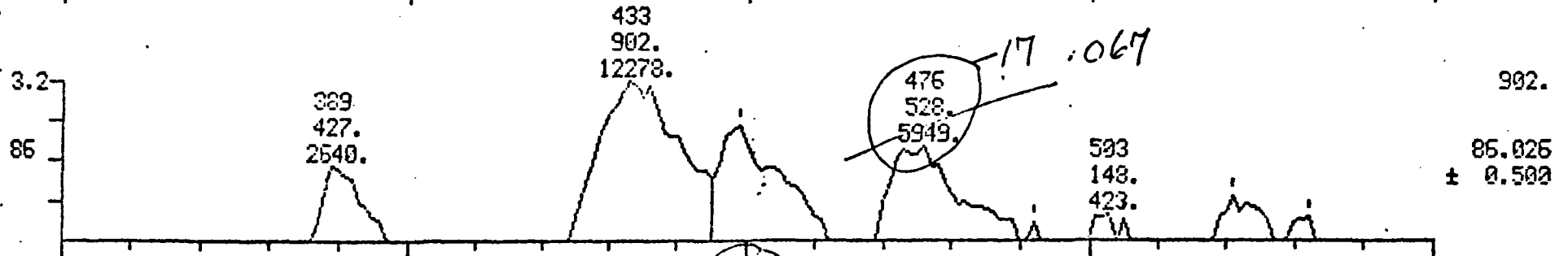
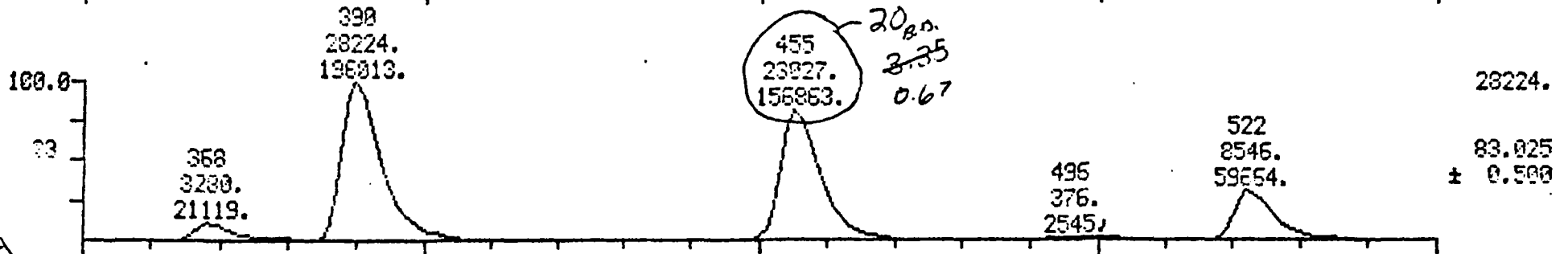
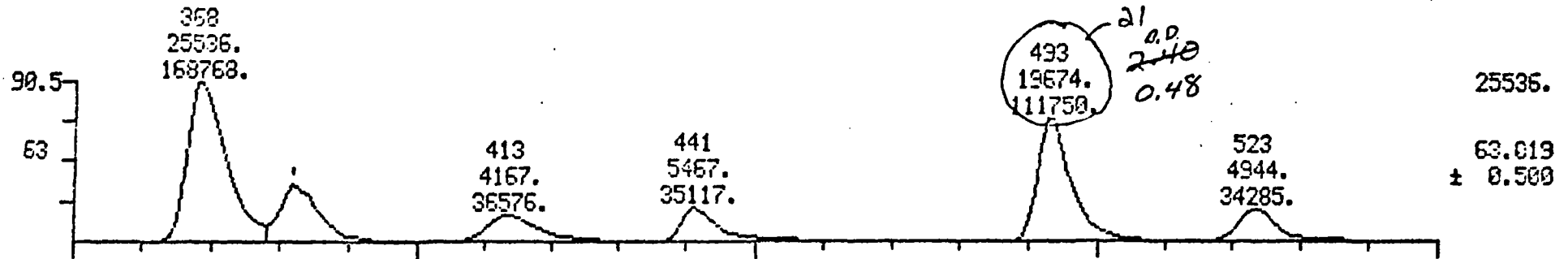
SCANS 300 TO 430



MASS CHROMATOGRAMS
 03/24/85 7:45:00
 SAMPLE: 58PF8 CASE 3931

DATA: U2716

SCANS 350 TO 550

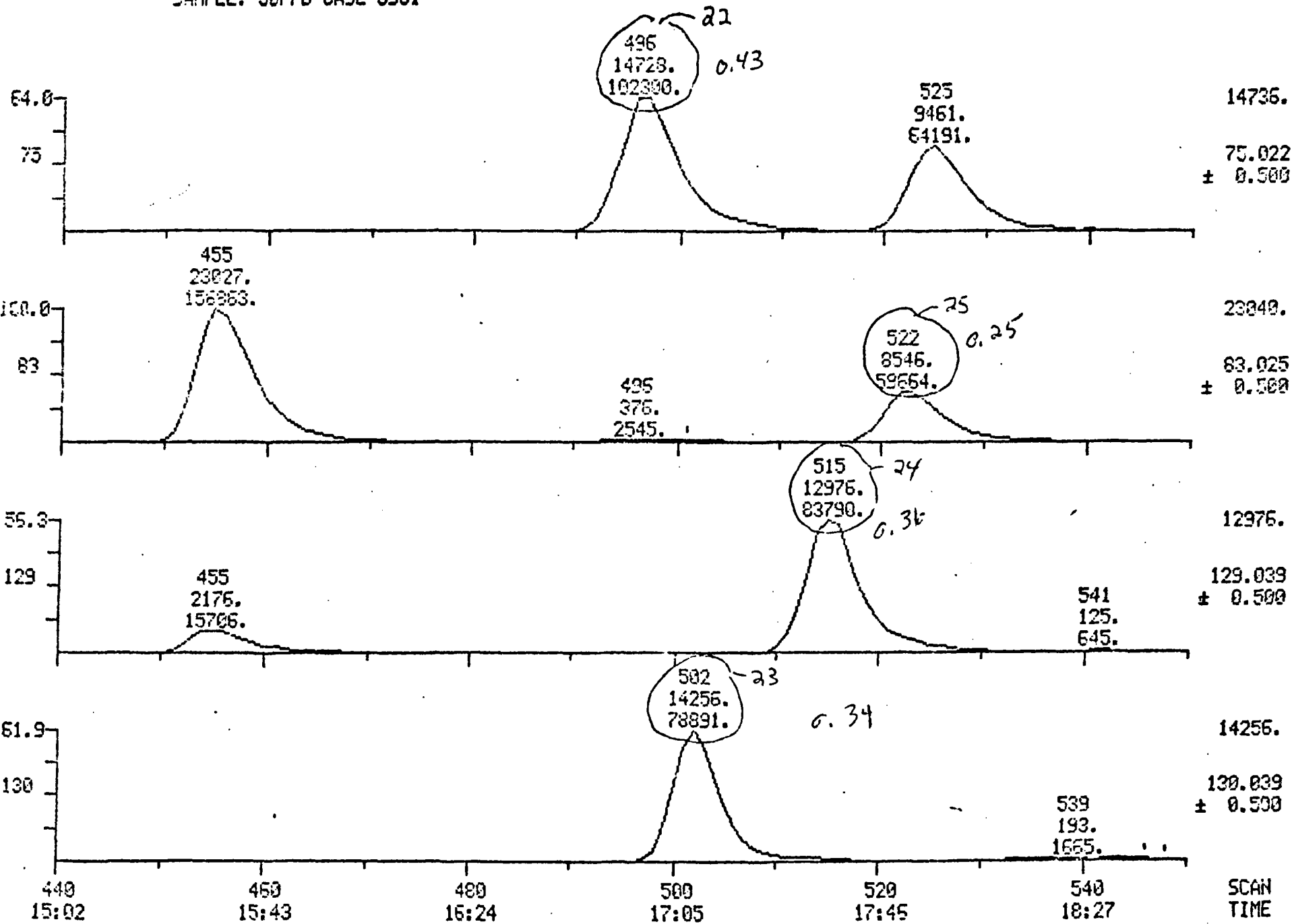


350 11:57 400 13:40 450 15:22 500 17:05 550 18:47 TIME

MASS CHROMATOGRAMS
 03/04/00 7:45:00
 SAMPLE: 50FFB CASE 3931

DATA: U2716

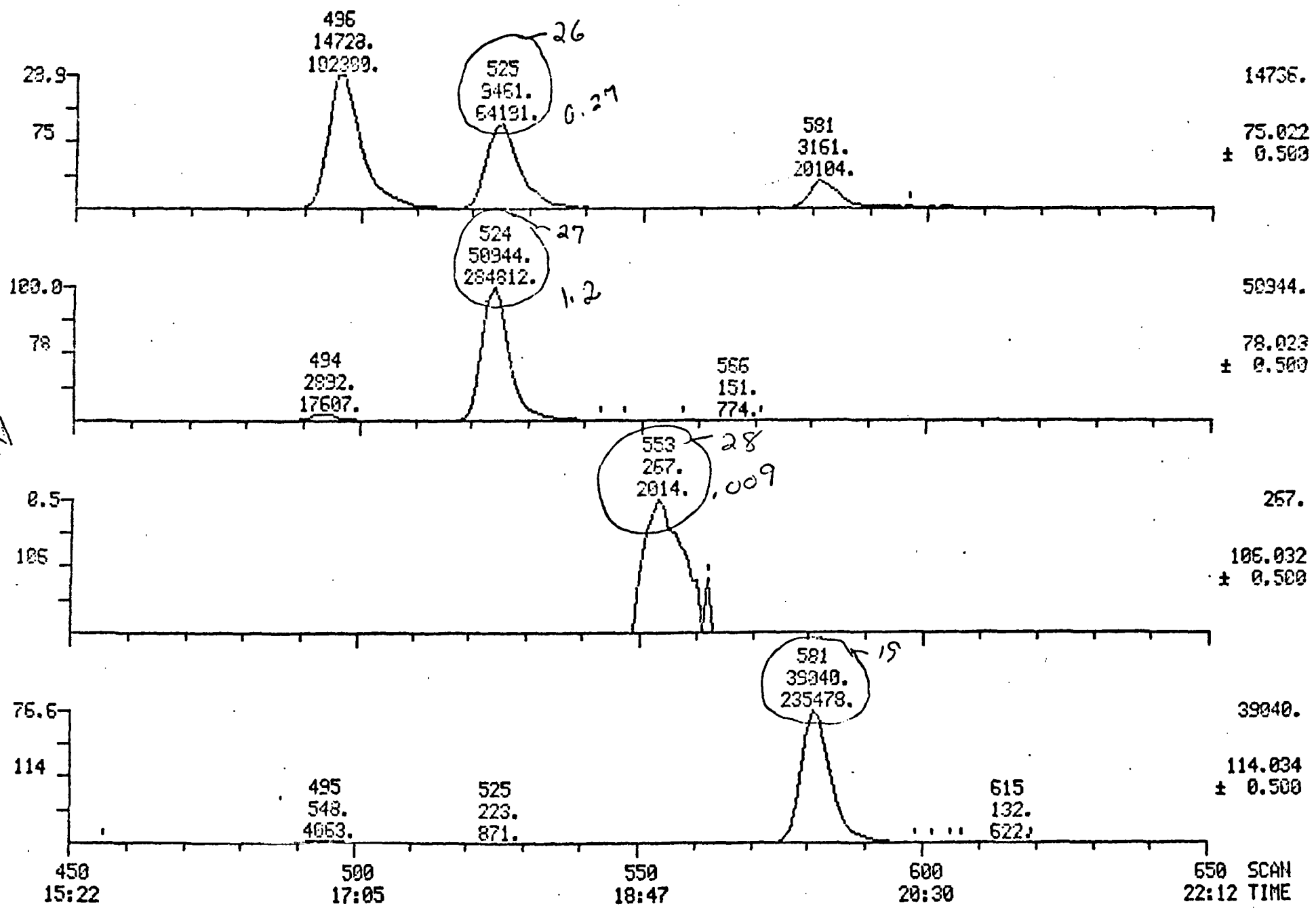
SCANS 440 TO 550



MASS CHROMATOGRAMS
 03/04/95 7:45:09
 SAMPLE: 50PPB CASE 3931

DATA: U2716

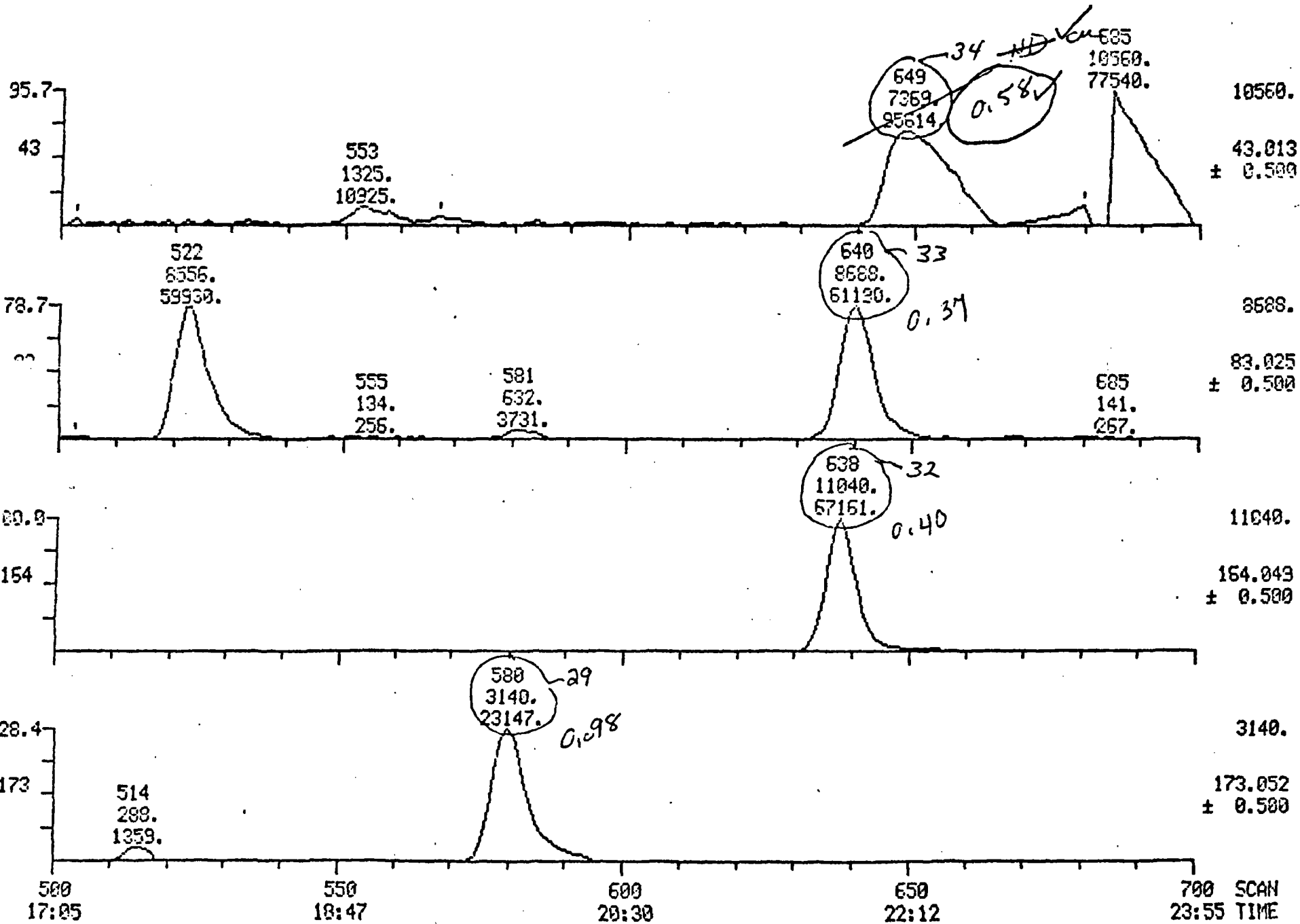
SCANS 450 TO 650



MASS CHROMATOGRAMS
 03/04/85 7:45:00
 SAMPLE: 50PPB CASE 3931

DATA: U2715

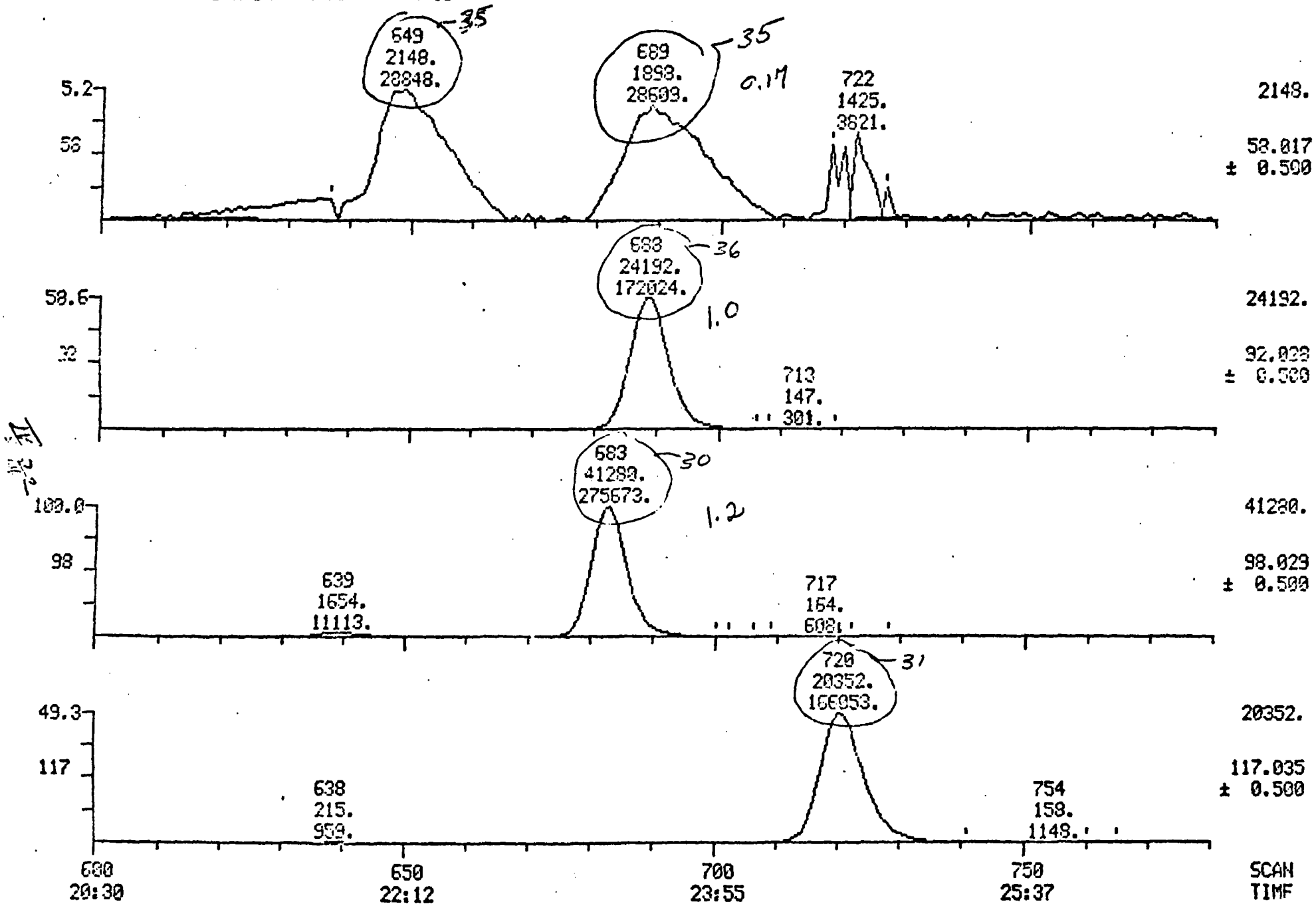
SCANS 500 TO 700



MASS CHROMATOGRAMS
03/04/95 7:45:00
SAMPLE: 50PPB CASE 3931

DATA: U2716

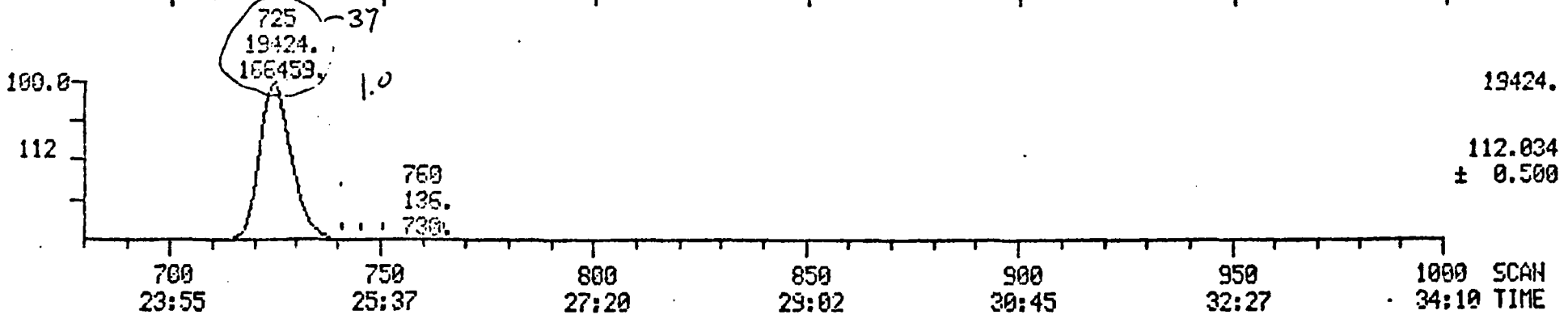
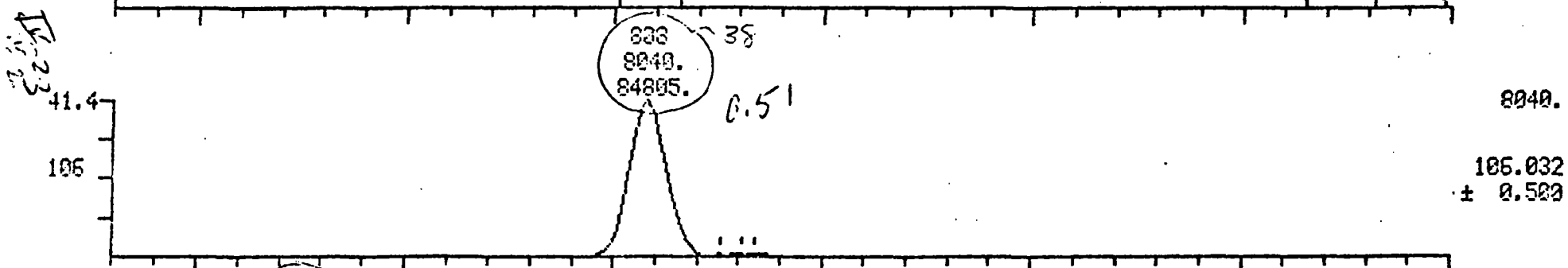
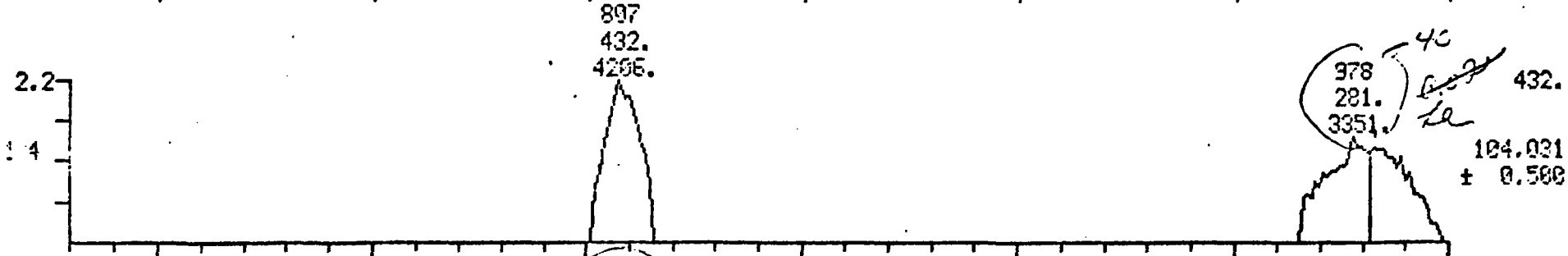
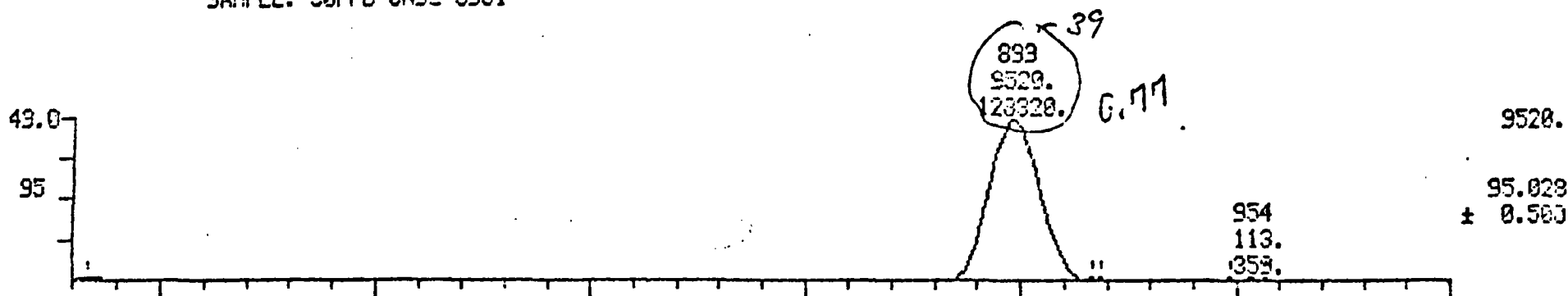
SCANS 600 TO 780



MASS CHROMATOGRAMS
03/04/65 7:45:00
SAMPLE: 50PPB CASE 3931

DATA: U2715

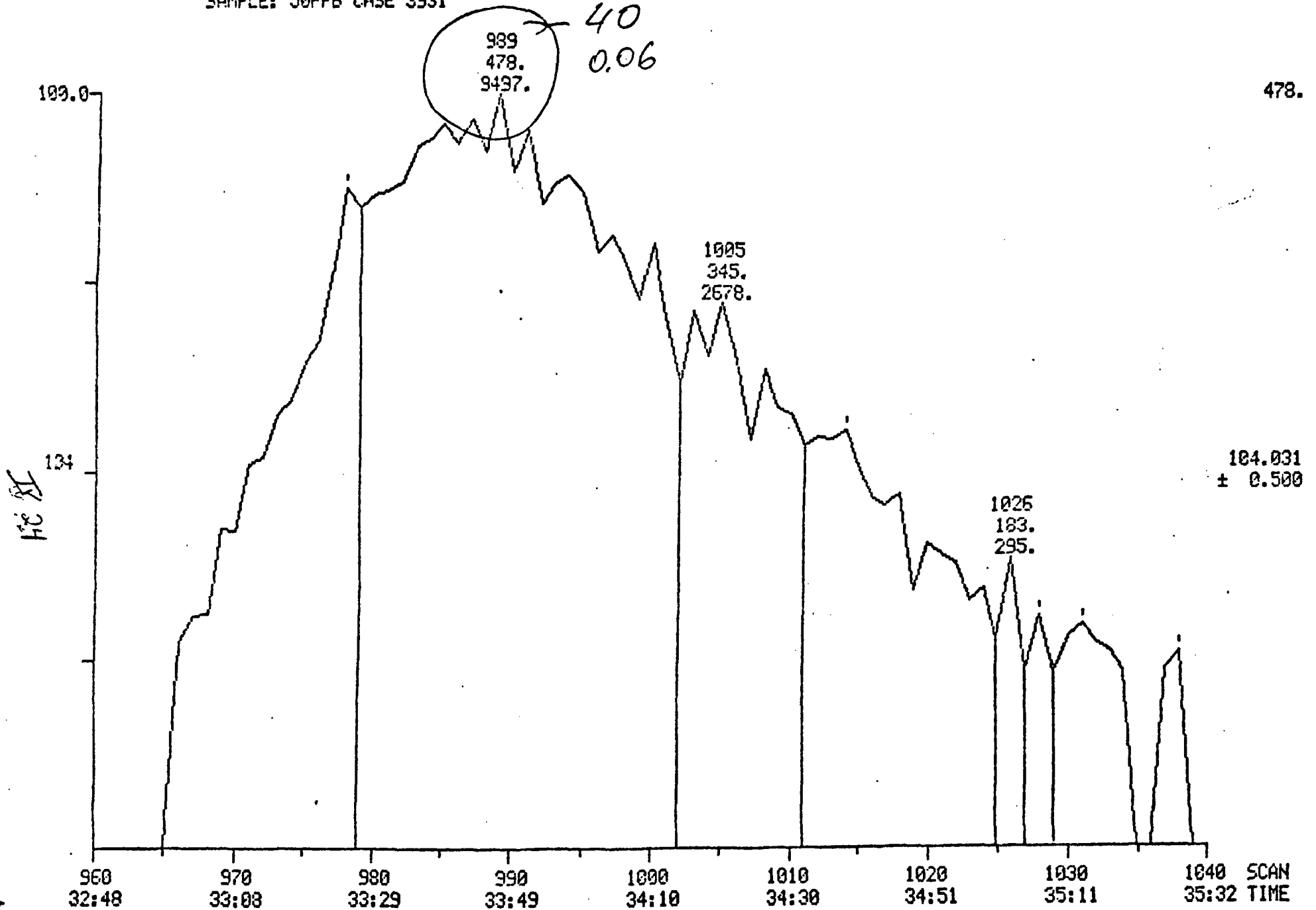
SCANS 680 TO 1000



MASS CHROMATOGRAM
03/04/25 7:45:09
SAMPLE: 50PPB CASE 3931

DATA: U2716

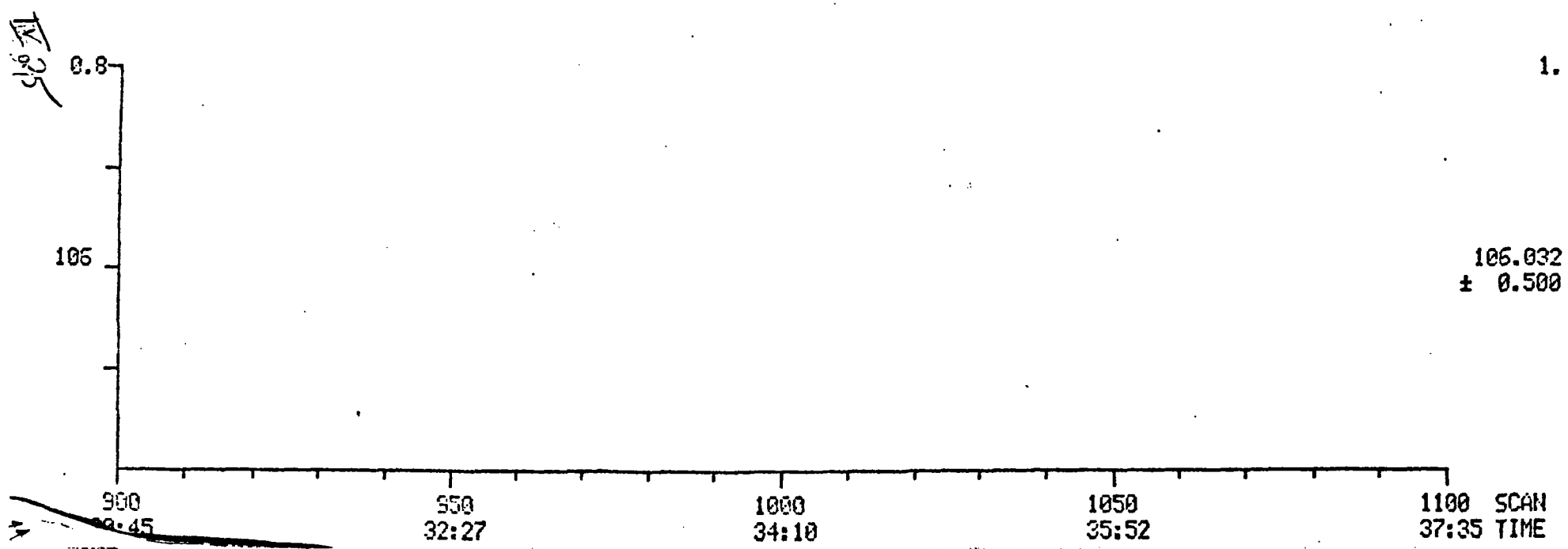
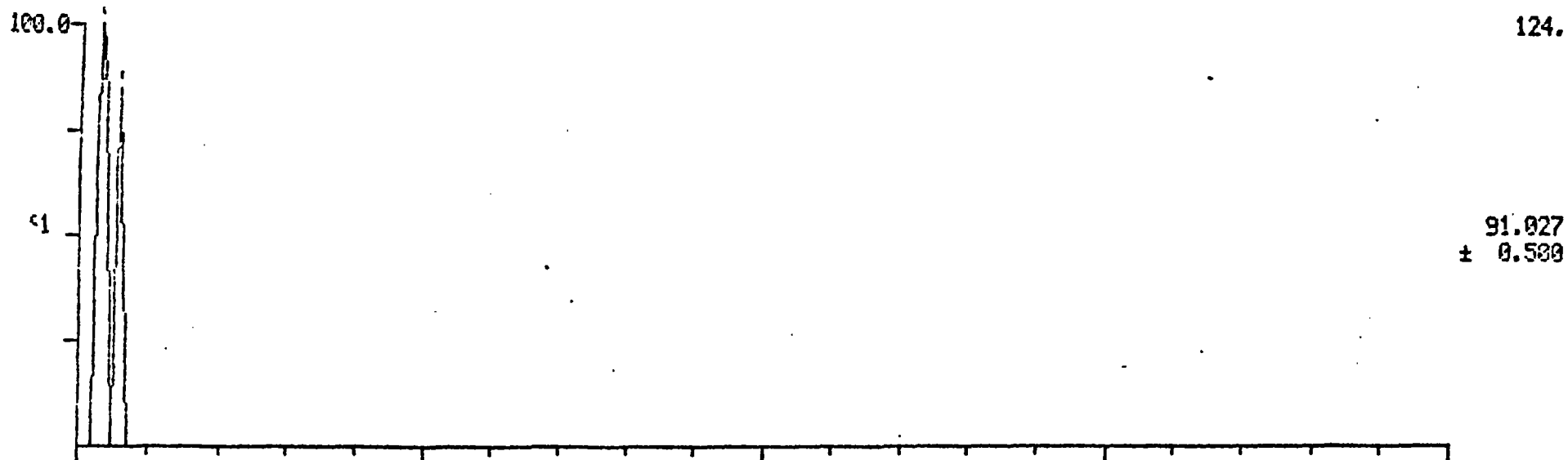
SCANS 960 TO 1040



MASS CHROMATOGRAMS
03/04/85 7:45:00
SAMPLE: 50PPB CASE 3331

DATA: U2716

SCANS 900 TO 1100



Contractor: GCA
 Contract No: 68-01-676-7
 Instrument ID: Fungus Mt. OWA

Time: 15:24
 Laboratory ID: 13743
 Initial Calibration Date: 3/3/85

Minimum RF for SPCC is 0.300 Maximum %D for CCC is 25%

Compound	RF	RF ₅₀	%D	CCC	SPCC
Chloromethane	1.2	1.1	8.3		..
Bromomethane	1.2	1.2	-0		
Vinyl Chloride	1.1	1.2	9.0	.	
Chloroethane	0.84	1.1	31		
Methylene Chloride	1.7	2.2	29		
Acetone	2.1	0.31	10		
Carbon Disulfide	3.3	3.5	6.1		
1,1-Dichloroethene	1.2	1.5	25	*	
1,1-Dichloroethane	3.2	4.0	25		..
Trans-1,2-Dichloroethene	1.6	1.9	19		
Chloroform	4.3	5.3	23	.	
1,2-Dichloroethane	4.0	4.4	10		
2-Butanone	1.2	0.34	133.26		
1,1,1-Trichloroethane	3.8	3.9	2.1		
Carbon Tetrachloride	2.6	2.1	19		
Vinyl Acetate	0.53	0.20	42		
Bromodichloromethane	0.94	0.83	12		
1,2-Dichloropropane	0.42	0.53	26	.	
Trans-1,3-Dichloropropene	0.47	0.53	13		
Trichloroethene	0.37	0.38	2.7		
Dibromochloromethane	0.56	0.40	28		
1,1,2-Trichloroethane	0.29	0.33	14		
Benzene	1.2	1.4	17		
cis-1,3-Dichloropropene	0.33	0.39	18		
2-Chloroethylvinyl ether	0.01	0.04	130		
Bromoform	0.28	0.10	14		..
2-Hexanone	0.13	0.21	47		
4-Methyl-2-Pentanone	0.78	0.68	14		
Tetrachloroethene	0.48	0.44	9.3		
1,1,2,2-Tetrachloroethane	0.38	0.48	26		..
Toluene	0.98	1.1	12	.	
Chlorobenzene	0.77	1.1	13		..
Ethylbenzene	0.47	0.57	21	.	
Styrene	0.27	0.06	78		
Total Xylenes	ND	ND*			

RF₅₀ - Response Factor from daily standard (5) at 50 ug/l
 RF - Average Response Factor from initial calibration Form VI

%D - Percent Difference
 CCC - Calibration Check Compounds (.)
 SPCC - System Performance Check Compounds (..)

Form VII + See narrative
 ND - See narrative.

JW 7 CM
JW 26

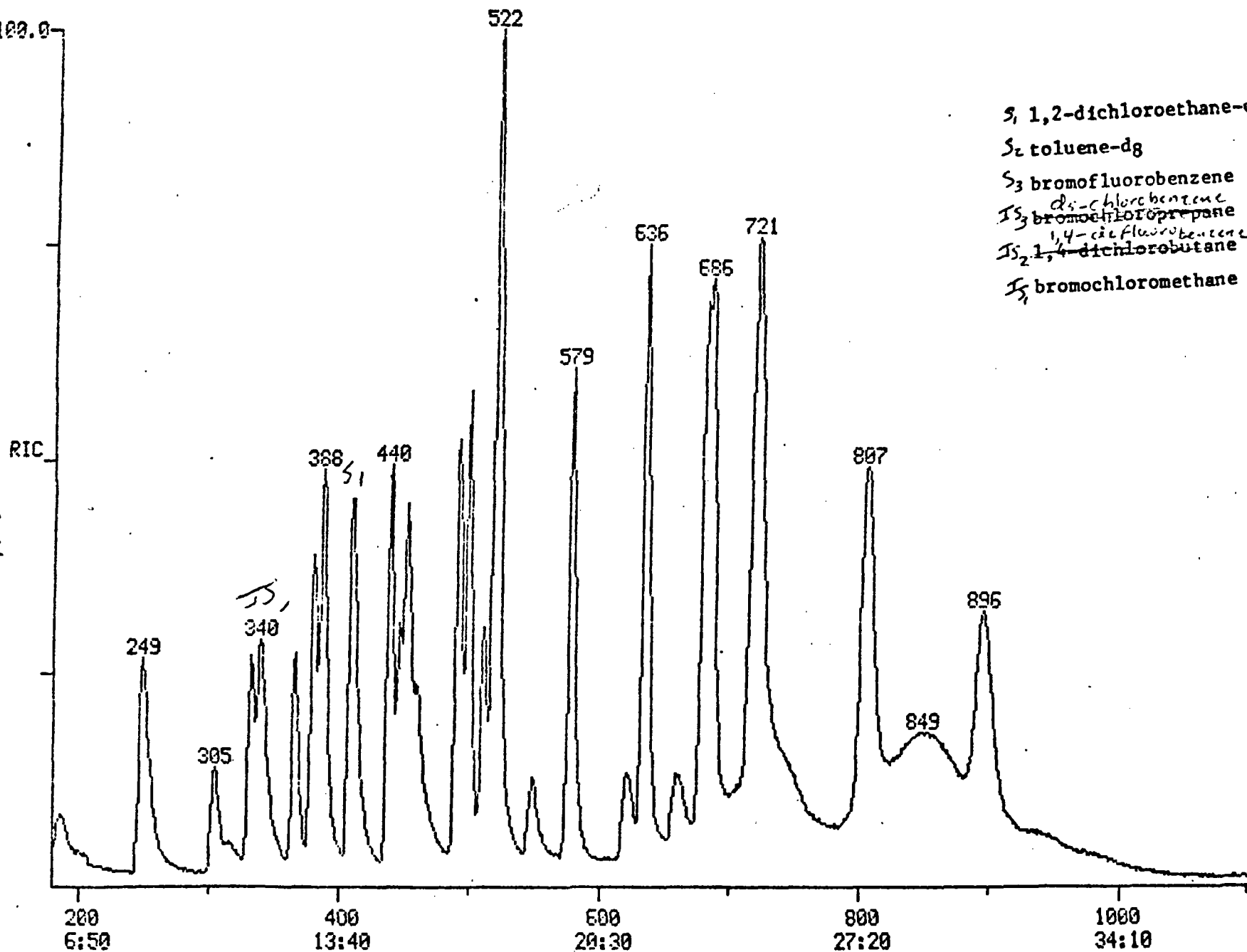
RIC
03/05/85 15:24:00
SAMPLE: 50 PPB STD CASE 3969

DATA: U2743

SCANS 100 TO 1100

156160.

116 II



- S₁ 1,2-dichloroethane-d₄
- S₂ toluene-d₈
- S₃ bromofluorobenzene
- S₄ ~~bromochloropropane~~
di-chlorobenzene
- S₅ ~~1,4-dichlorobutane~~
1,4-difluorobenzene
- S₆ bromochloromethane

SCAN
TIME

SILICAN ORGANICS III HEATER ANALYZER
CALCULATION REPORT FILE: V2190

AS: V2190.TE
08/09 10:28:00

FILE:
SUBMITTED BY: ANALYST:

UNIT=AREA(HIGHT) * REF. AREA / (REF. AREA(HIGHT) * RESP. FACT)
SP. FAC. FROM LIBRARY ENTRY

- 1 NAME
- 2 BROMOCHLOROMETHANE (INTERNAL STANDARD 1)
- 3 CHLOROMETHANE
- 4 BROMOMETHANE
- 5 METHYL CHLORIDE
- 6 CHLOROETHANE
- 7 NITRYLNE CHLORIDE
- 8 ACETONE
- 9 CARBON DISULFIDE
- 10 1,1-DICHLOROETHANE
- 11 1,2-DICHLOROETHANE
- 12 TRANS-1,2-DICHLOROETHENE
- 13 CHLOROFORM
- 14 D1-1,2-DICHLOROETHANE (SURROGATE 1)
- 15 1,2-DICHLOROETHENE
- 16 2-BUTANONE
- 17 1,1,1-TRICHLOROETHANE
- 18 CARBON TETRACHLORIDE
- 19 VINYL ACETATE
- 20 BROMODICHLOROMETHANE
- 21 1,2-DICHLOROPROPANE
- 22 TRANS-1,2-DICHLOROPROPENE
- 23 TRICHLOROETHENE
- 24 DIBROMODICHLOROMETHANE
- 25 1,1,2-TRICHLOROETHANE
- 26 BENZENE
- 27 CIS-1,2-DICHLOROPROPENE
- 28 2-CHLOROETHYL METHYL ETHER
- 29 1,4-DIFLUOROBENZENE (INTERNAL STANDARD 2)
- 30 BROMOFORM
- 31 4-METHYLPENTANONE
- 32 1,1,2,2-TETRACHLOROETHANE
- 33 TETRACHLOROETHENE
- 34 2-HEXANONE
- 35 D2-1,2-DICHLOROETHANE (SURROGATE 2)
- 36 TOLUENE
- 37 D3-1,2-DICHLOROETHANE (INTERNAL STANDARD 3)
- 38 CHLOROBENZENE
- 39 ETHYL BENZENE
- 40 4-FLUOROBENZENE (SURROGATE 3)
- 41 STYRENE
- 42 TOTAL XYLENES

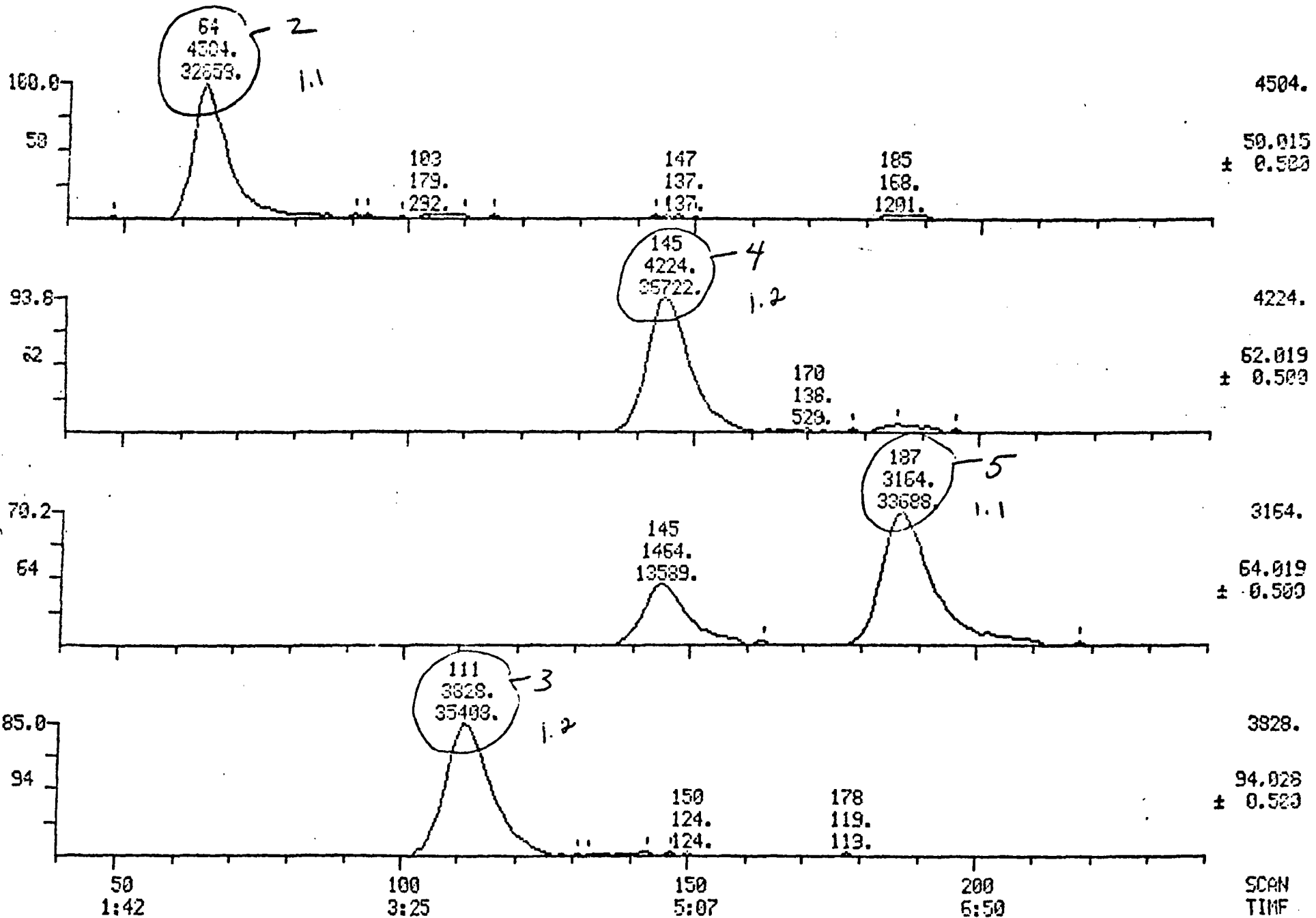
Handwritten signature

Handwritten number 28

MASS CHROMATOGRAMS
03/05/85 15:24:00
SAMPLE: 50 PFB STD CASE 3963

DATA: U2743

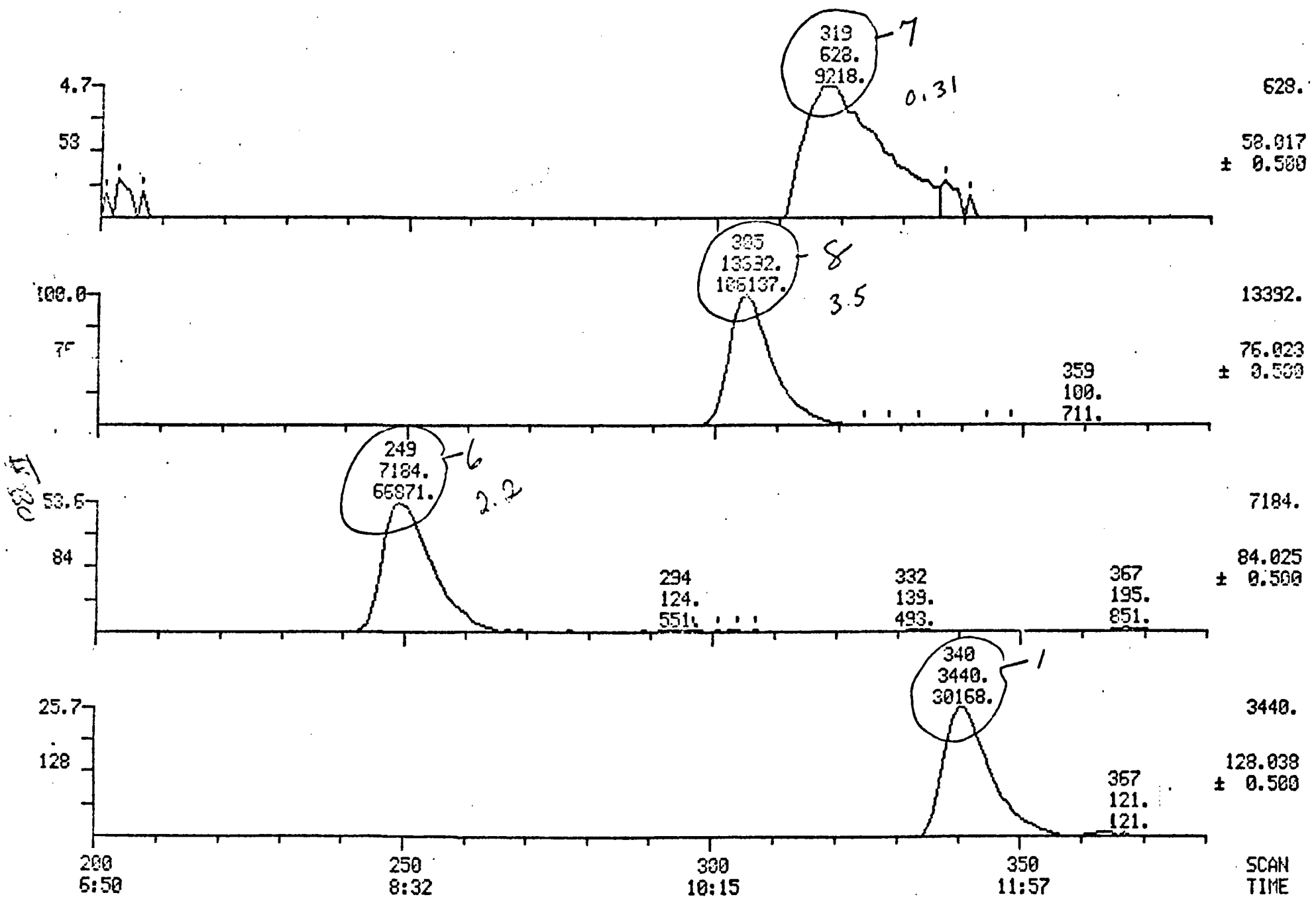
SCANS 40 TO 240



MASS CHROMATOGRAMS
03/05/95 15:24:00
SAMPLE: 50 PPB STD CASE 3969

DATA: U2743.

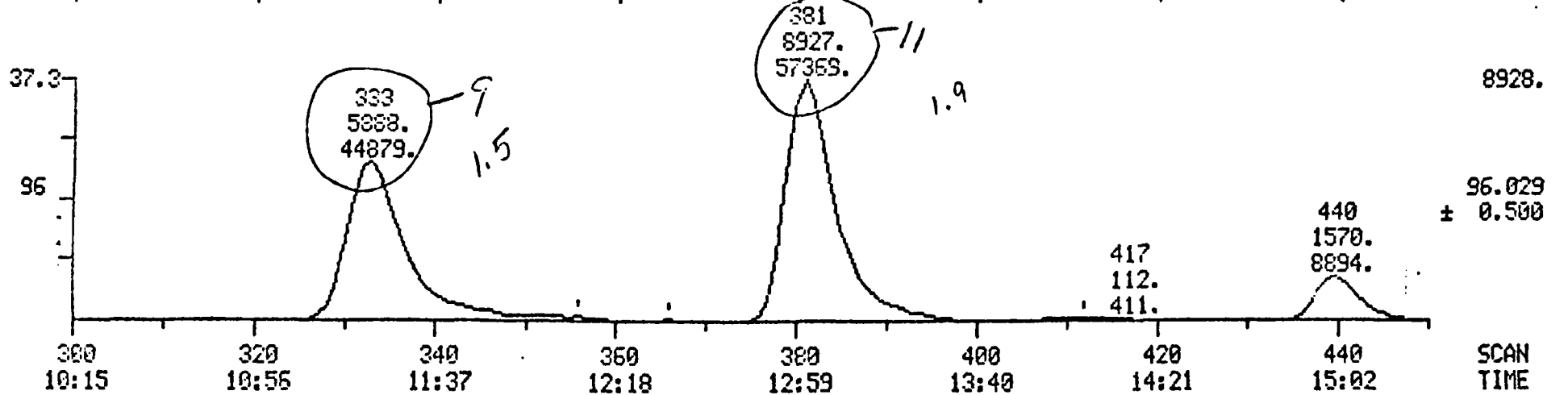
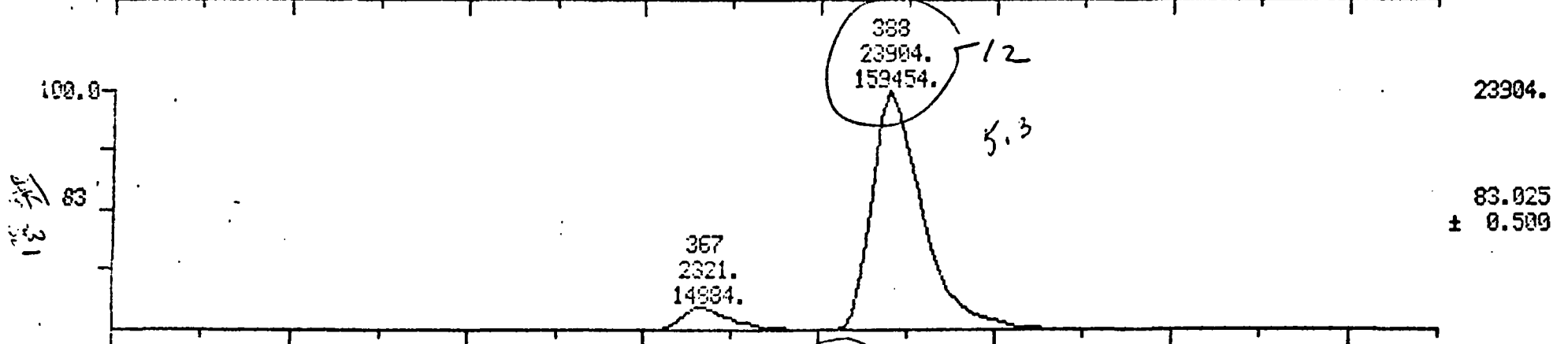
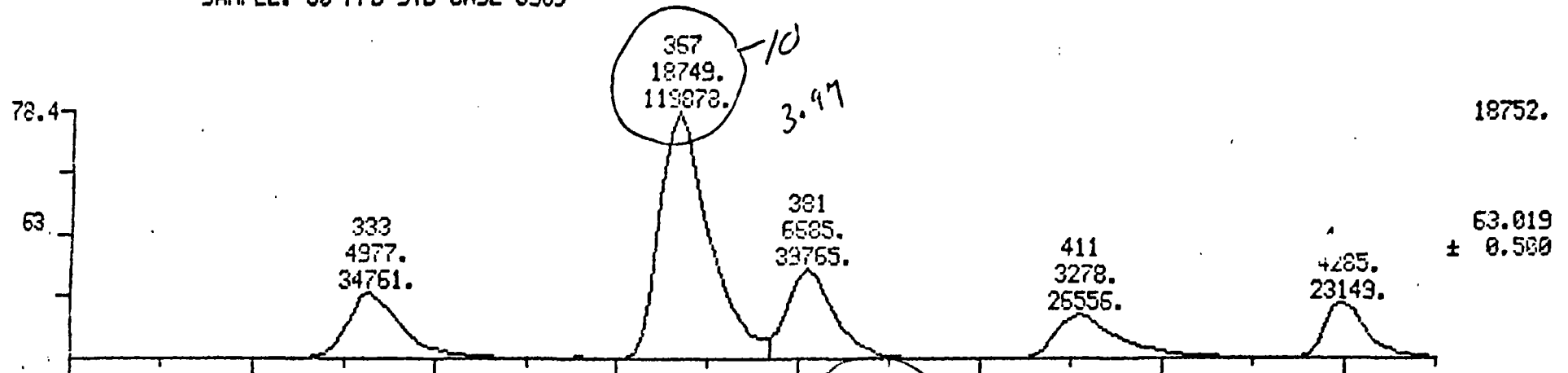
SCANS 200 TO 380



MASS CHROMATOGRAMS
03/05/85 15:24:00
SAMPLE: 50 FPB STD CASE 3969

DATA: U2743

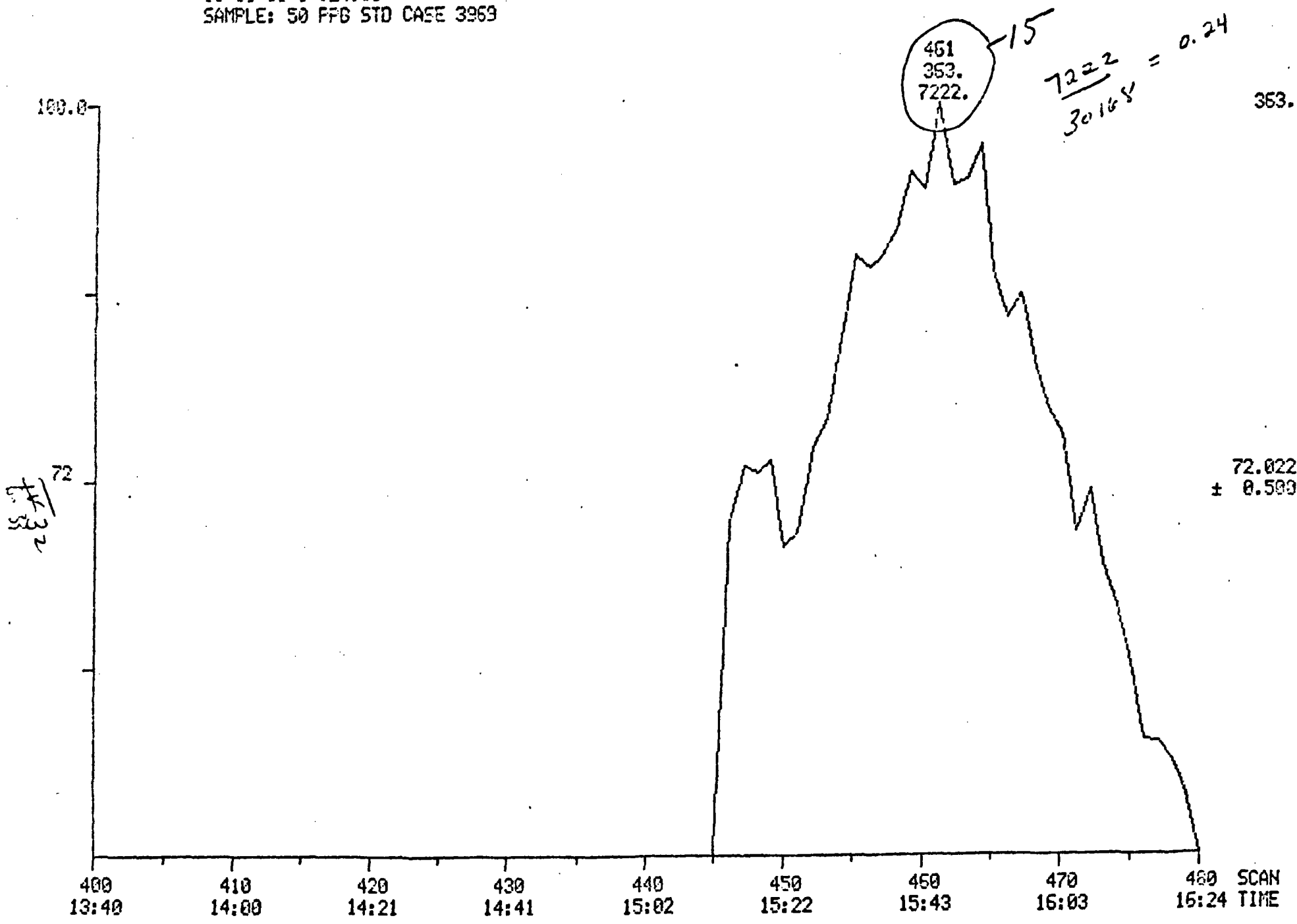
SCANS 300 TO 450



MASS CHROMATOGRAM
03/05/65 15:24:00
SAMPLE: 50 FFB STD CASE 3959

DATA: U2743

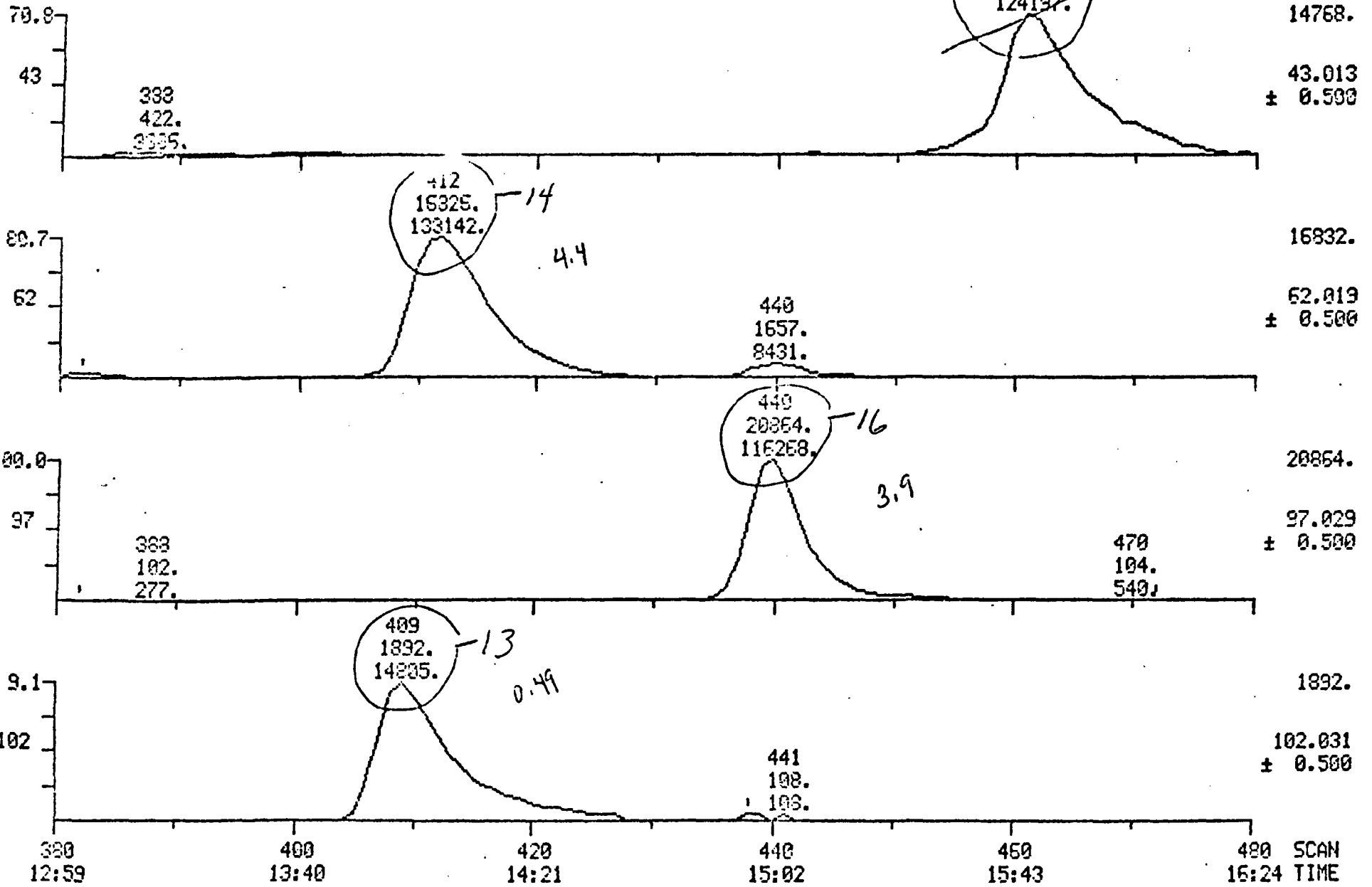
SCANS 400 TO 480



MASS CHROMATOGRAMS
 03/05/85 15:24:03
 SAMPLE: 50 PPB STD CASE 3959

DATA: U2743

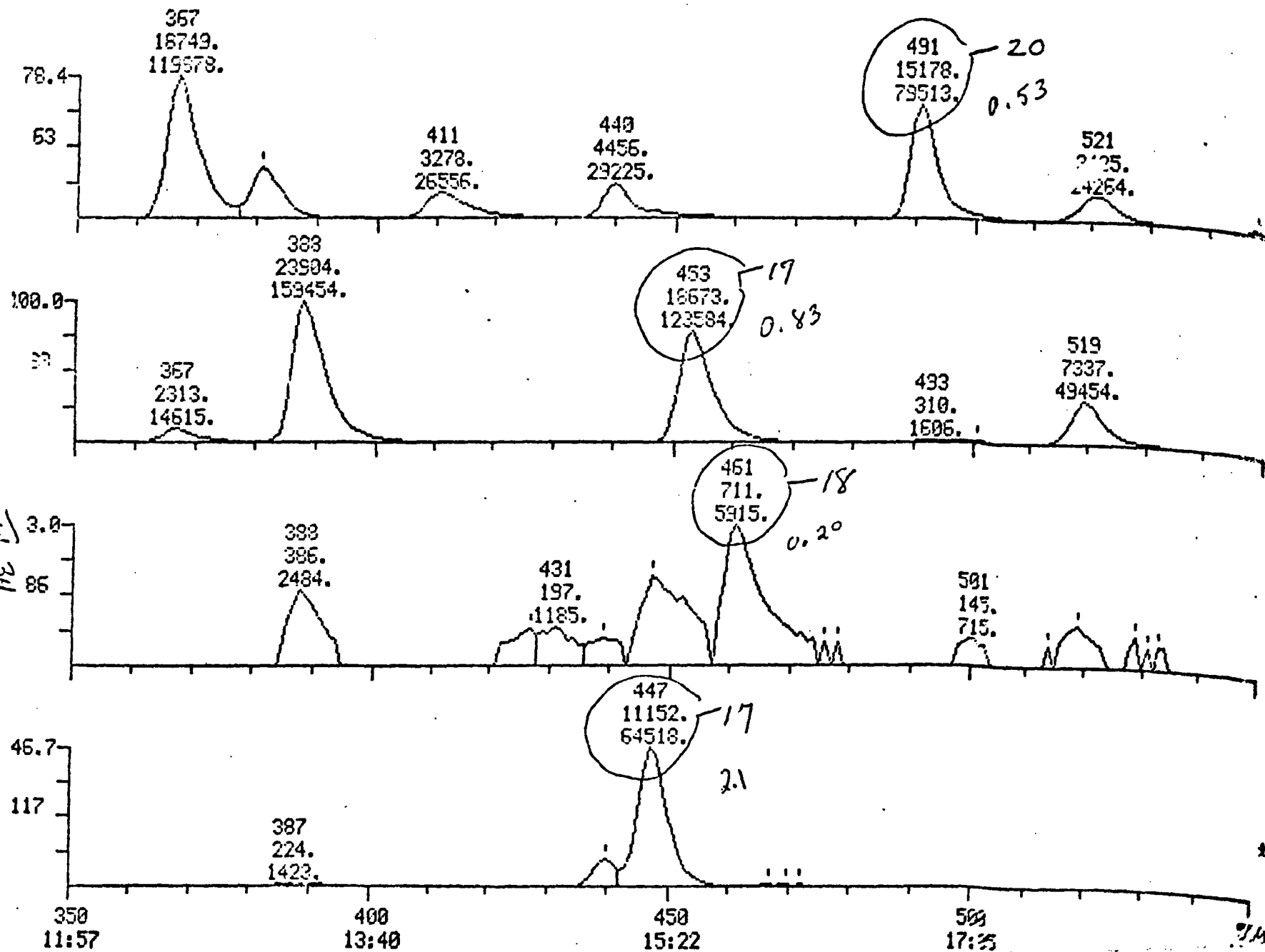
SCANS 330 TO 450



MASS CHROMATOGRAMS
03/05/65 15:24:00
SAMPLE: 50 PFB STD CASE 3959

DATA: U2743

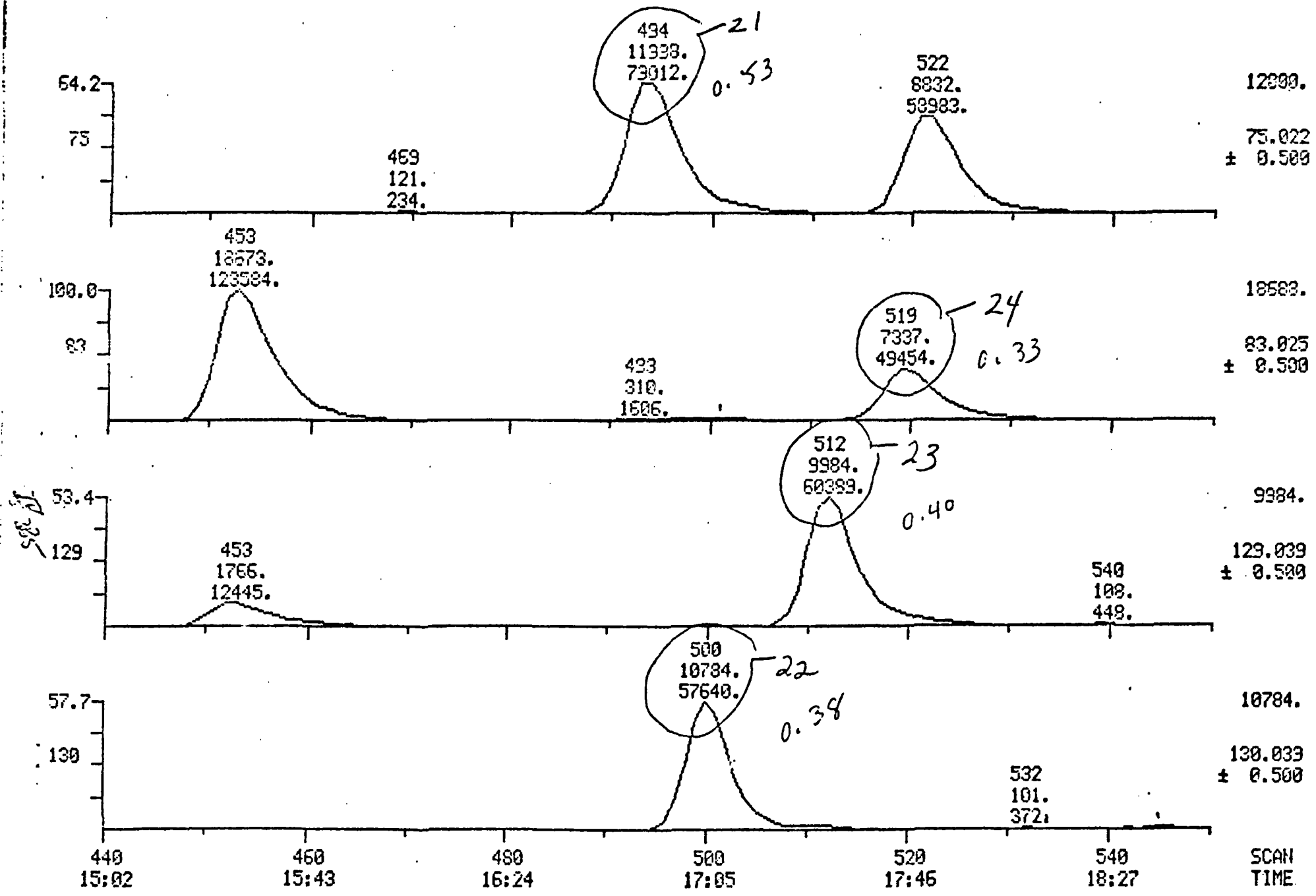
SCANS 352 10



MASS CHROMATOGRAMS
 03/05/85 15:24:00
 SAMPLE: 50 FPB STD CASE 3969

DATA: U2743

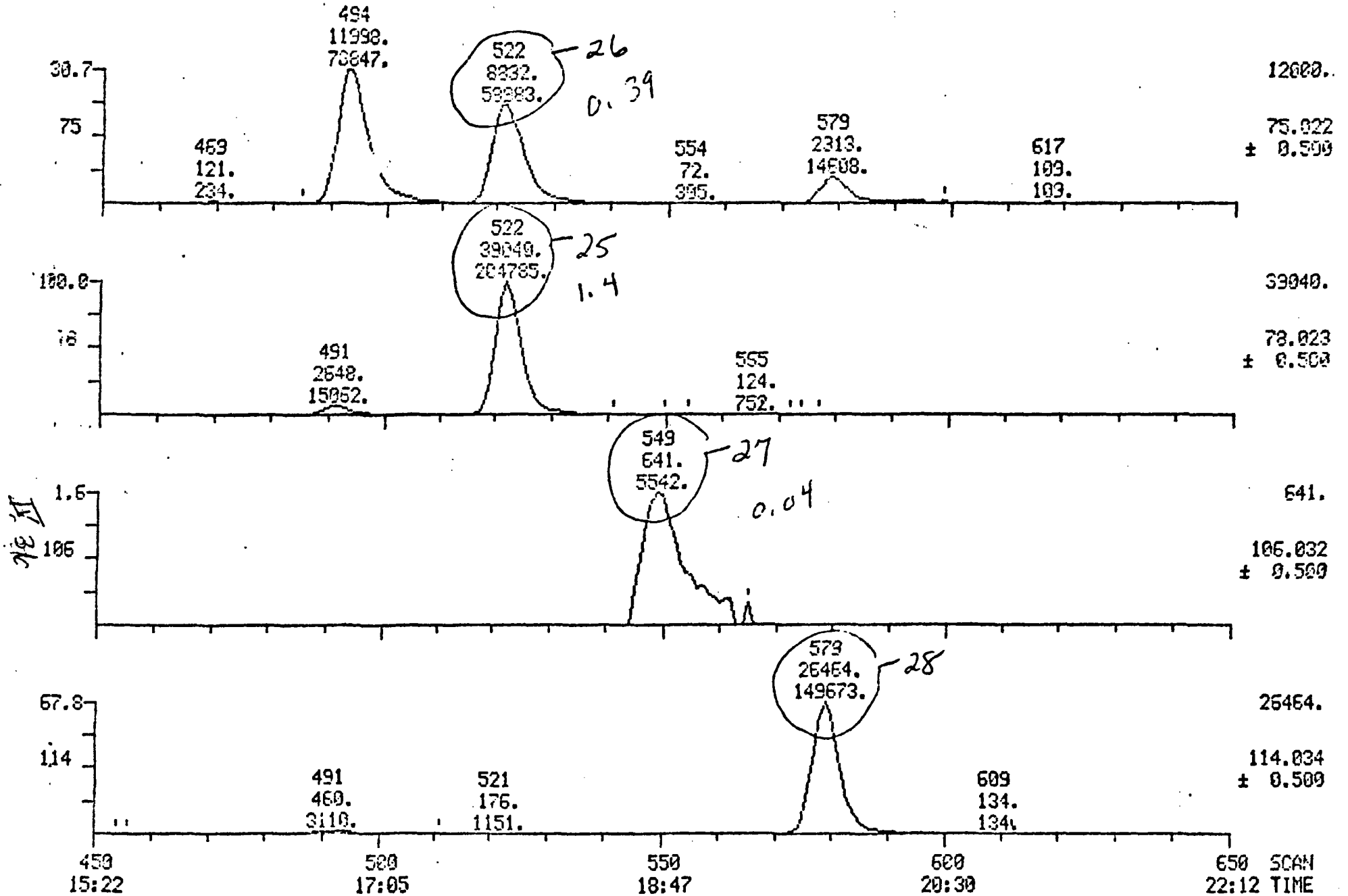
SCANS 440 TO 550



MASS CHROMATOGRAMS
03/05/85 15:24:00
SAMPLE: 50 PPB STD CASE 3969

DATA: U2743

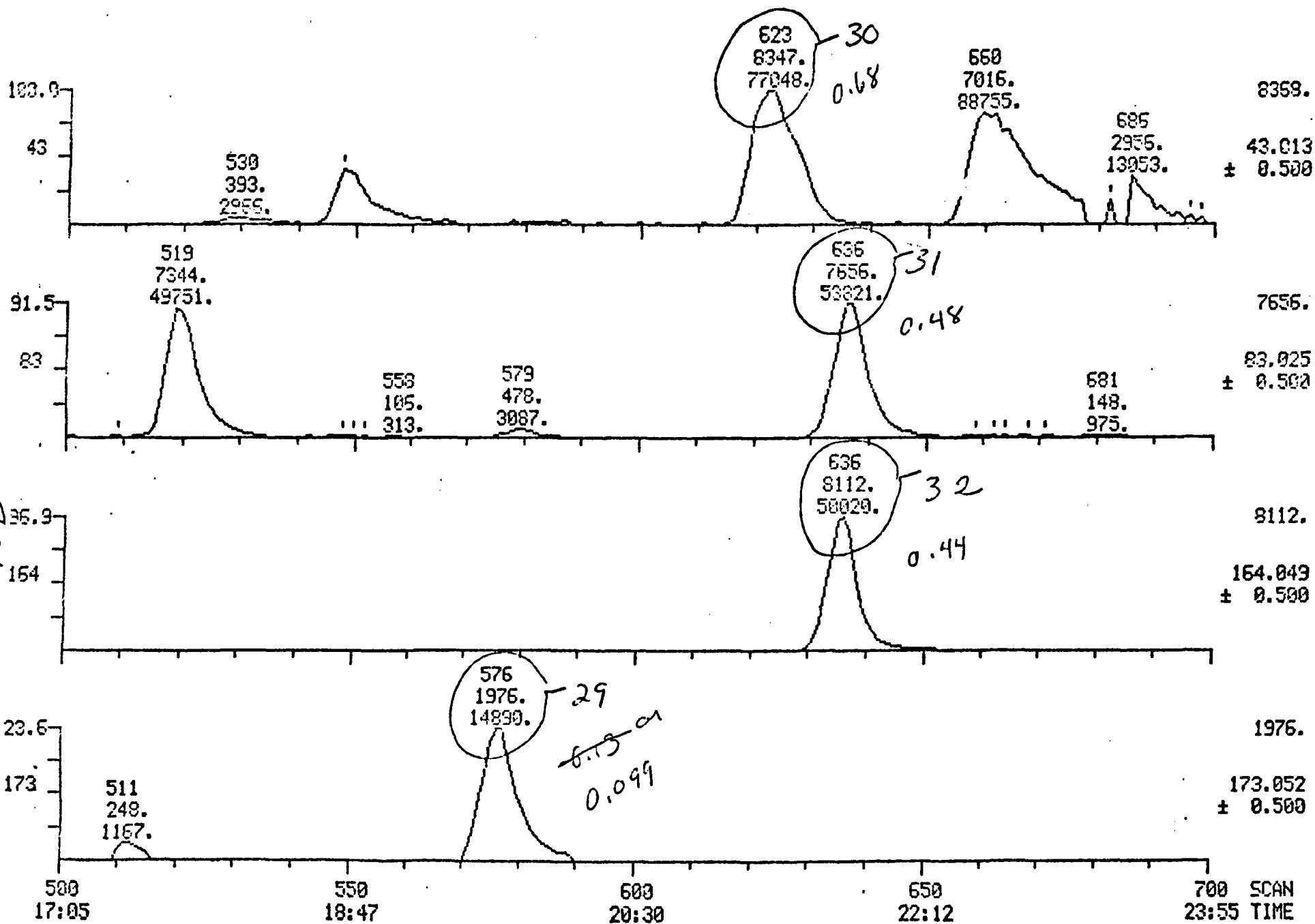
SCANS 450 TO 550



MASS CHROMATOGRAMS
03/25/85 15:24:00
SAMPLE: 50 PFB STD CASE 3969

DATA: U2743

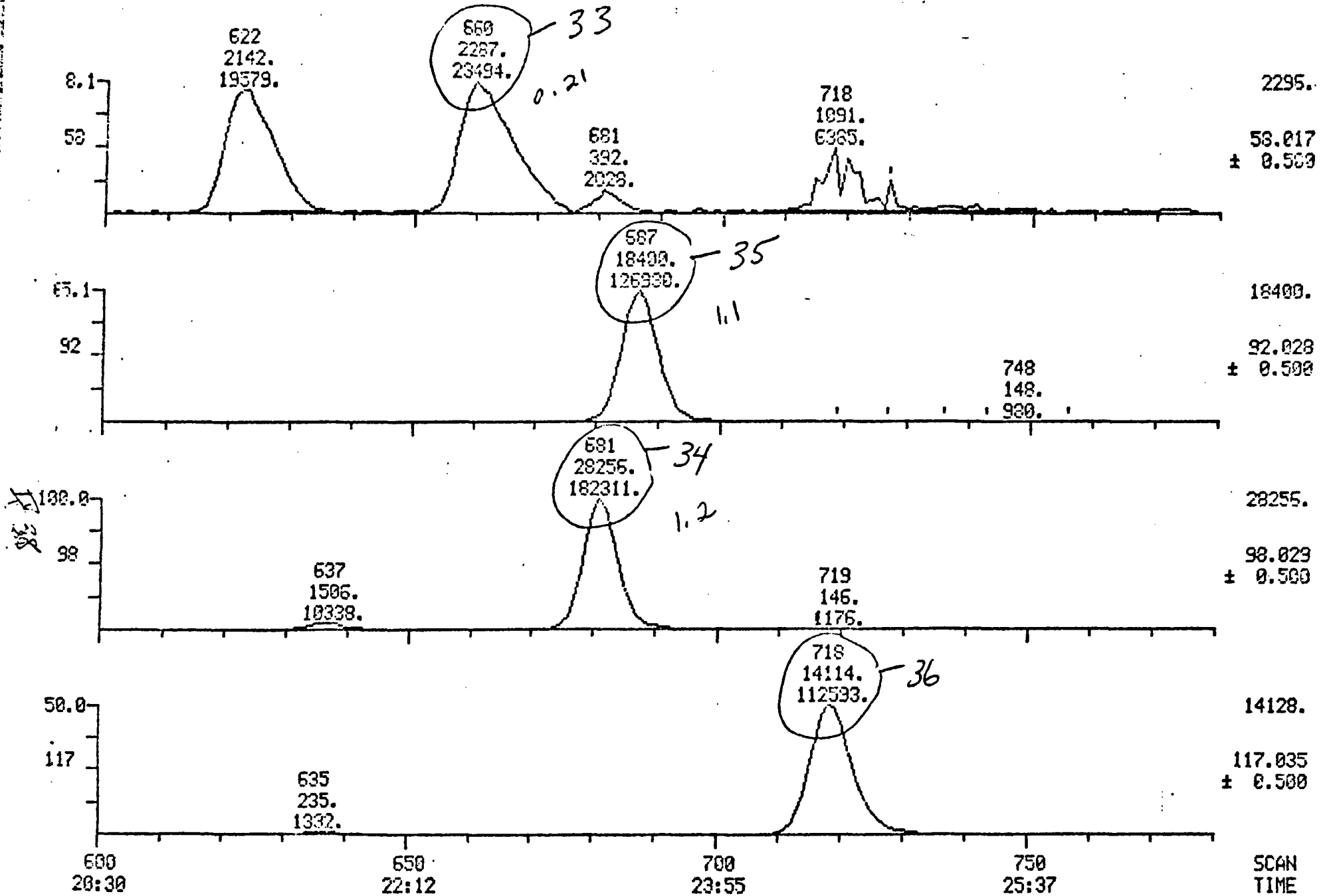
SCANS 500 TO 700



MASS CHROMATOGRAMS
03/05/85 15:24:00
SAMPLE: 50 PPB STD CASE 3959

DATA: U2743

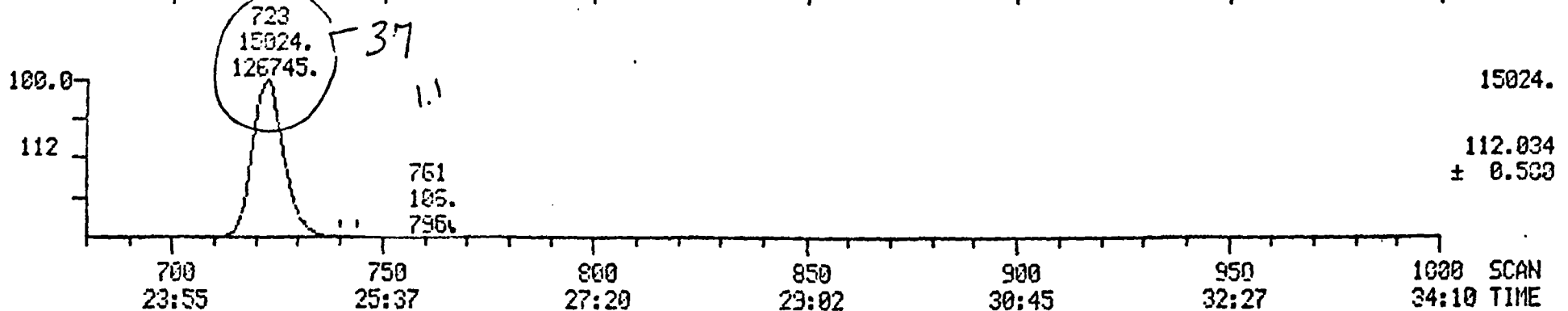
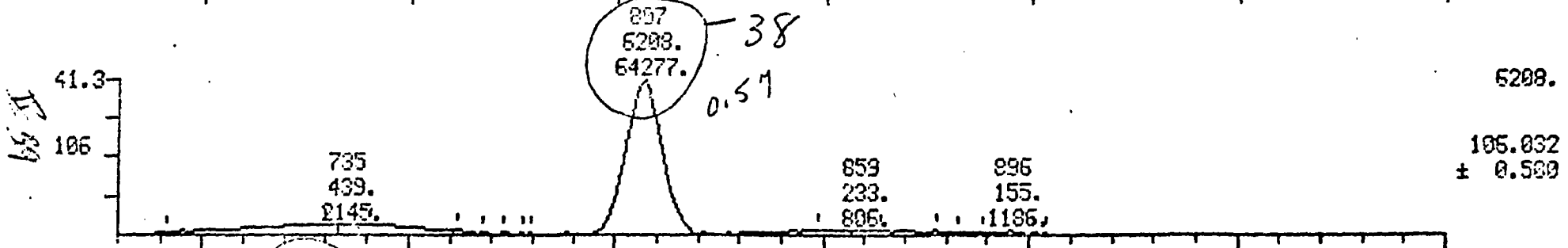
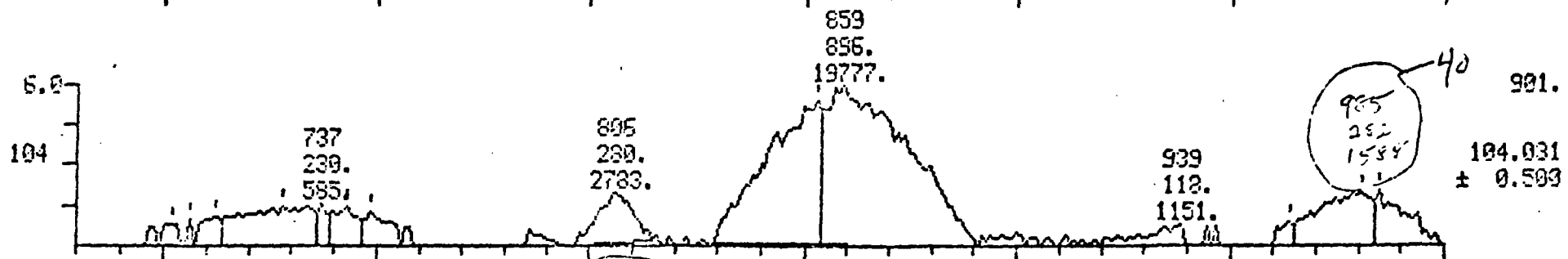
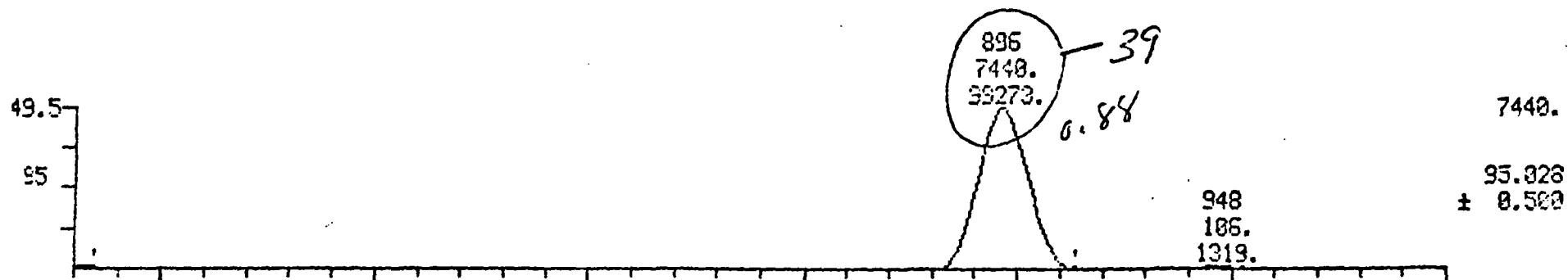
SCANS 600 TO 720



MASS CHROMATOGRAMS
03/05/05 15:24:03
SAMPLE: 50 PF8 STD CASE 3969

DATA: U2743

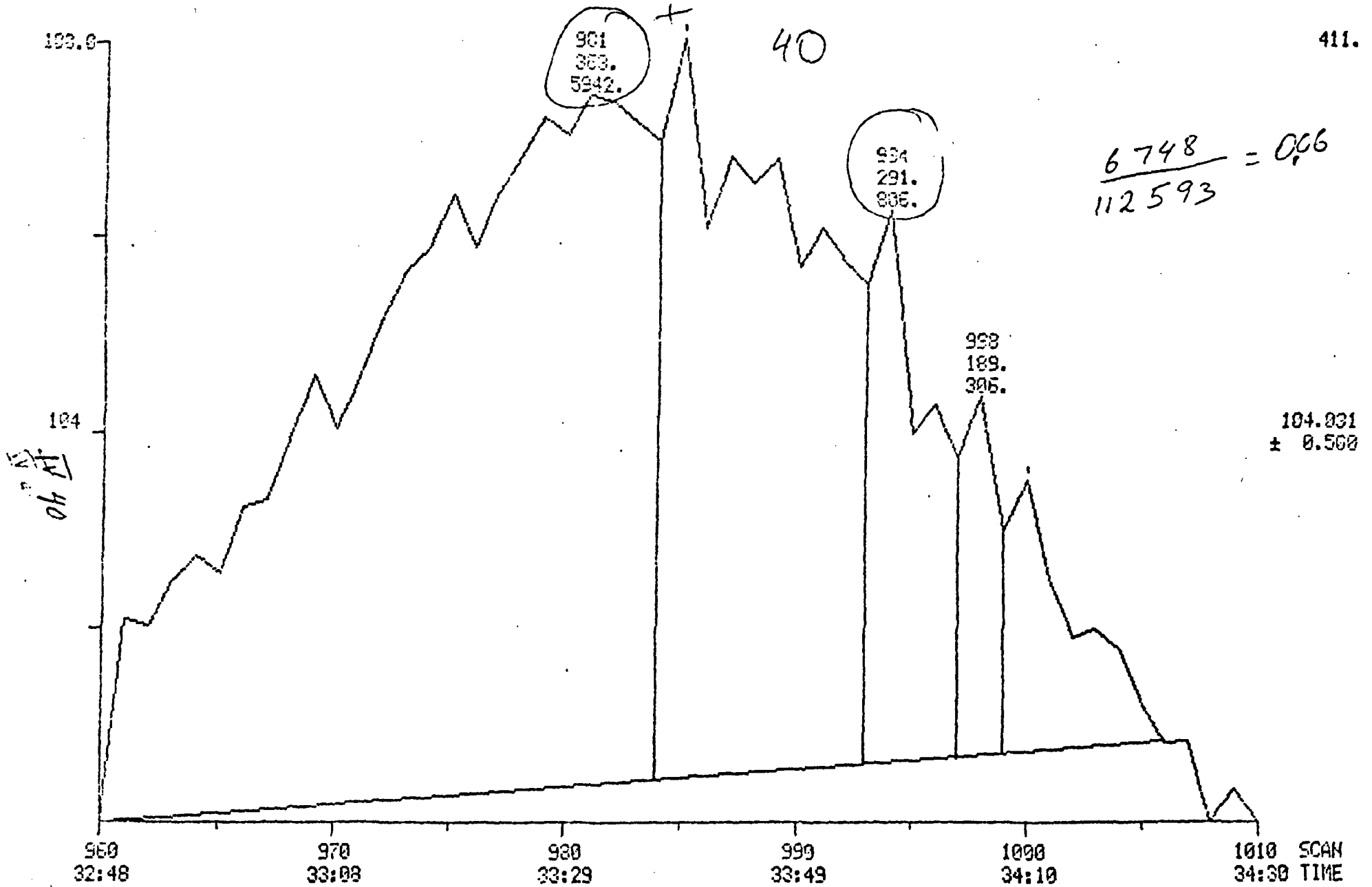
SCANS 680 TO 1000



MASS CHROMATOGRAM
03/05/85 15:24:00
SAMPLE: 50 PFB STD CASE 3369

DATA: U2743

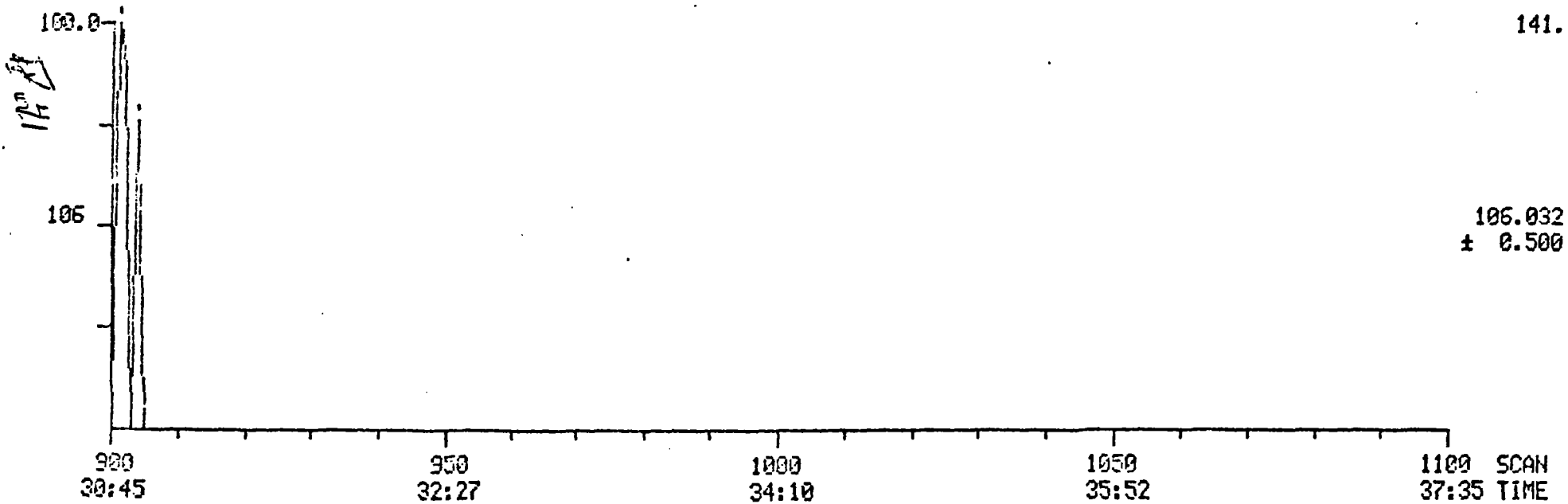
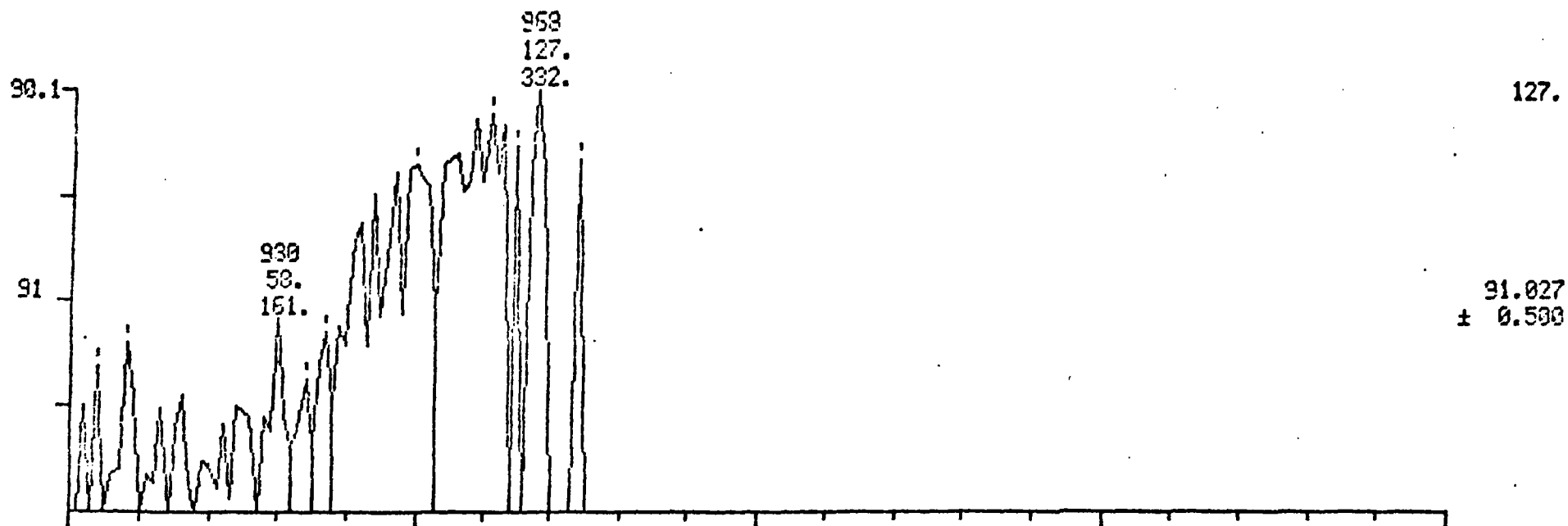
SCANS 959 TO 1019



MASS CHROMATOGRAMS
03/05/85 15:24:09
SAMPLE: 50 PPB STD CASE 3369

DATA: U2743

SCANS 900 TO 1100



Case No: 3969
 Contractor: GCA
 Contract No: 68-01-6767
 Instrument ID: Finnigan MAT OWA

Calibration Date: 3/6/85
 Time: 11:43
 Laboratory ID: V2755
 Initial Calibration Date: 3/3/85

Minimum RF for SPCC is 0.300 Maximum %D for CCC is 25%

Compound	RF	RF ₅₀	%D	CCC	SPCC
Chloromethane	1.2	0.88	27		..
Bromomethane	1.2	1.1	8.3		
Vinyl Chloride	1.1	1.0	9.1	*	
Chloroethane	1.84	0.93	11		
Methylene Chloride	1.7	2.4	41		
Acetone	1.21	0.24	14		
Carbon Disulfide	3.3	3.2	3.0		
1, 1-Dichloroethene	1.2	1.3	10	*	
1, 1-Dibromoethene	3.2	4.0	25		..
Trans-1, 2-Dichloroethene	1.6	1.9	19		
Chloroform	4.3	5.1	19	*	
1, 2-Dichloroethane	4.0	4.1	2.5		
2-Butanone	1.2	0.10	92		
1, 1, 1-Trichloroethane	3.8	3.5	7.9		
Carbon Tetrachloride	2.6	1.8	31		
Vinyl Acetate	0.13	0.15	15		
Bromochloromethane	0.94	0.77	18		
1, 2-Dichloropropane	0.42	0.52	24	*	
Trans-1, 3-Dichloropropene	0.47	0.49	2.0		
Trichloroethene	0.37	0.36	2.7		
Dibromochloromethane	0.54	0.35	35		
1, 1, 2-Trichloroethane	0.29	0.30	3.4		
Benzene	1.2	1.3	8.3		
cis-1, 3-Dichloropropene	0.33	0.32	3.0		
2-Chloroethylvinylether	0.01	0.03	200		
Bromoforn	0.28	0.08	74		..
2-Pentanone	0.13	0.19	46		
4-Methyl-2-Pentanone	0.78	0.64	18		
Tetrachloroethene	0.48	0.42	13		
1, 1, 2, 2-Tetrachloroethane	0.38	0.44	16		..
Toluene	0.98	1.1	12	*	
Chlorobenzene	0.97	1.1	13		..
Ethylbenzene	0.47	0.57	21	*	
Styrene	0.27	0.23	15		
Total Xylenes	ND	ND	ND		

RF₅₀ - Response Factor from daily standard (6) at 50 ug/l
 RF - Average Response Factor from impical calibration Form VI

%D - Percent Difference
 CCC - Calibration Check Compounds (*)
 SPCC - System Performance Check Compounds (..)

Form VII

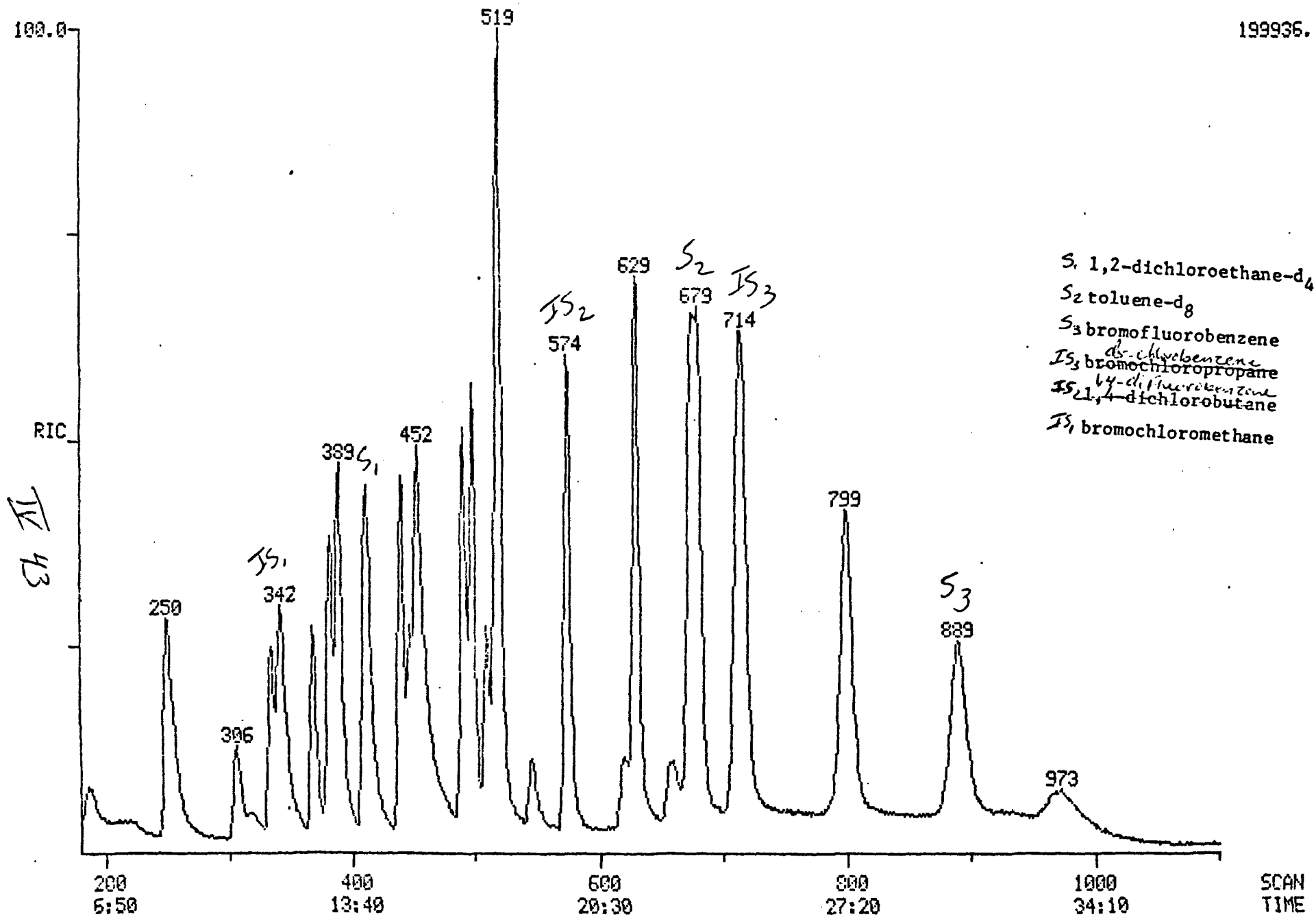
IV 42

RIC
03/05/85 11:43:00
SAMPLE: 50 PPB STD 3/6/85 Case 3969

DATA: U2755

SCANS 180 TO 1100

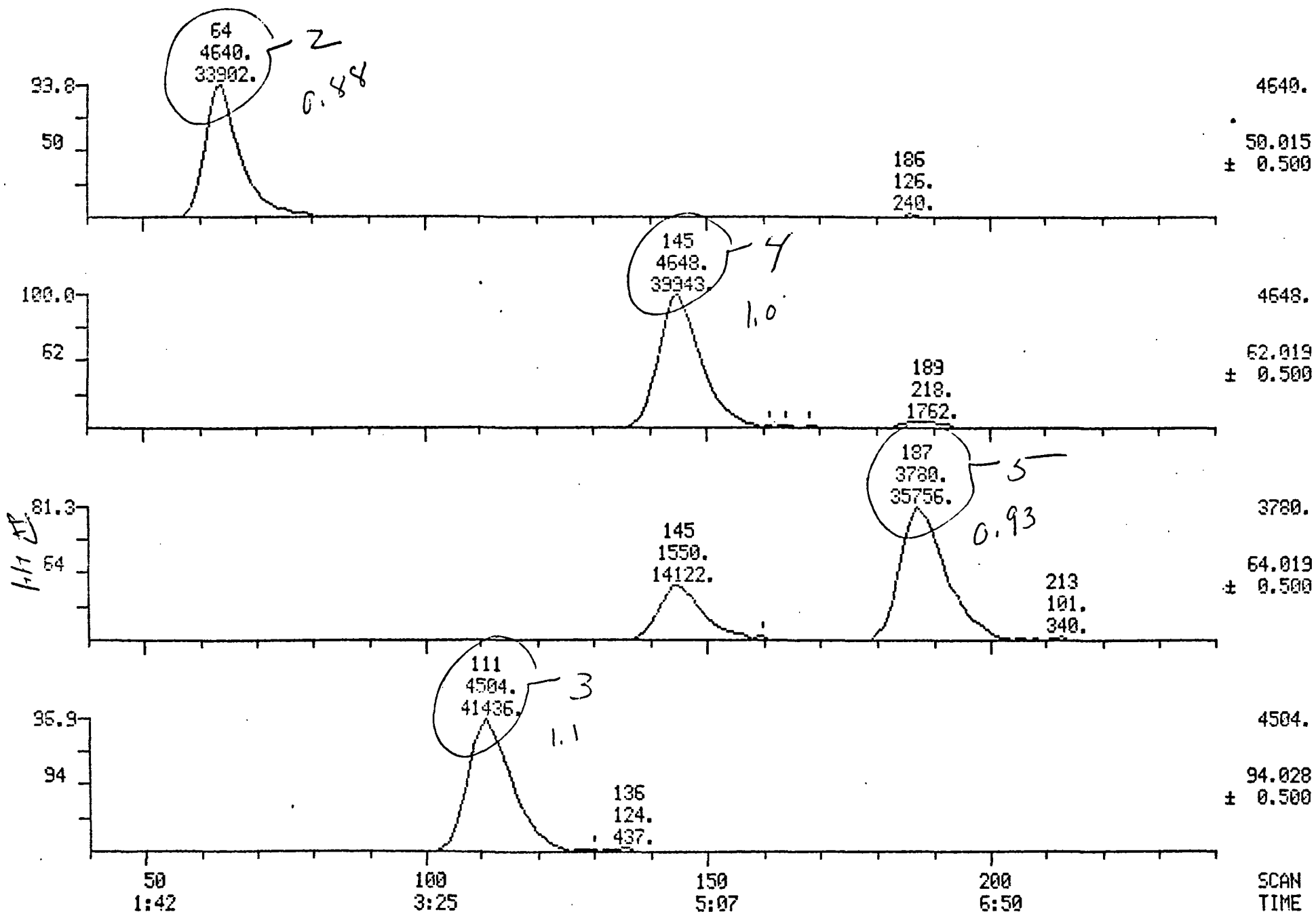
199936.



MASS CHROMATOGRAMS
03/06/85 11:43:00
SAMPLE: 50 PPB STD 3/6/85 Case 3969

DATA: U2755

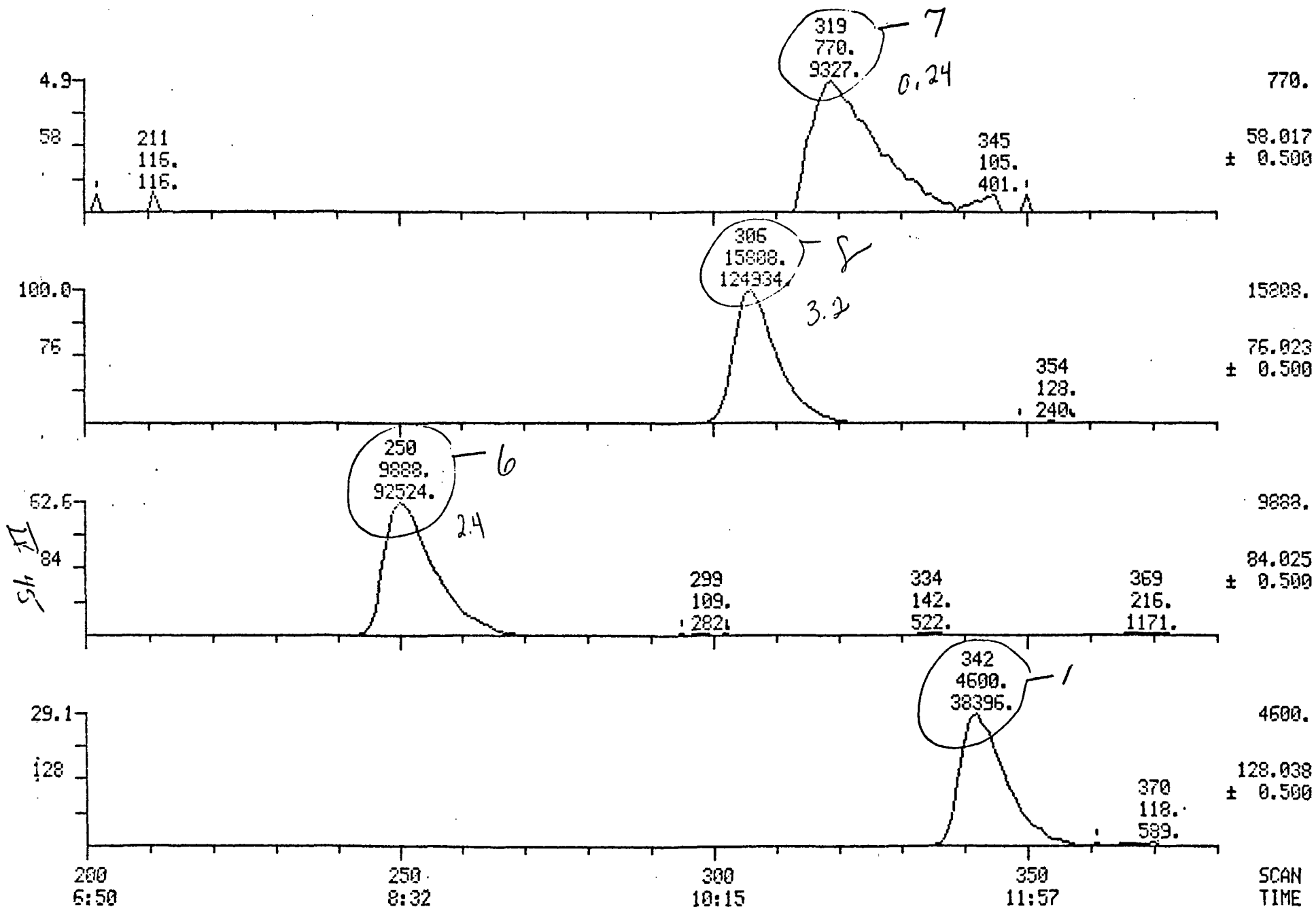
SCANS 40 TO 240



MASS CHROMATOGRAMS
 03/06/85 11:43:00
 SAMPLE: 50 PFB STD 3/6/85

DATA: U2755

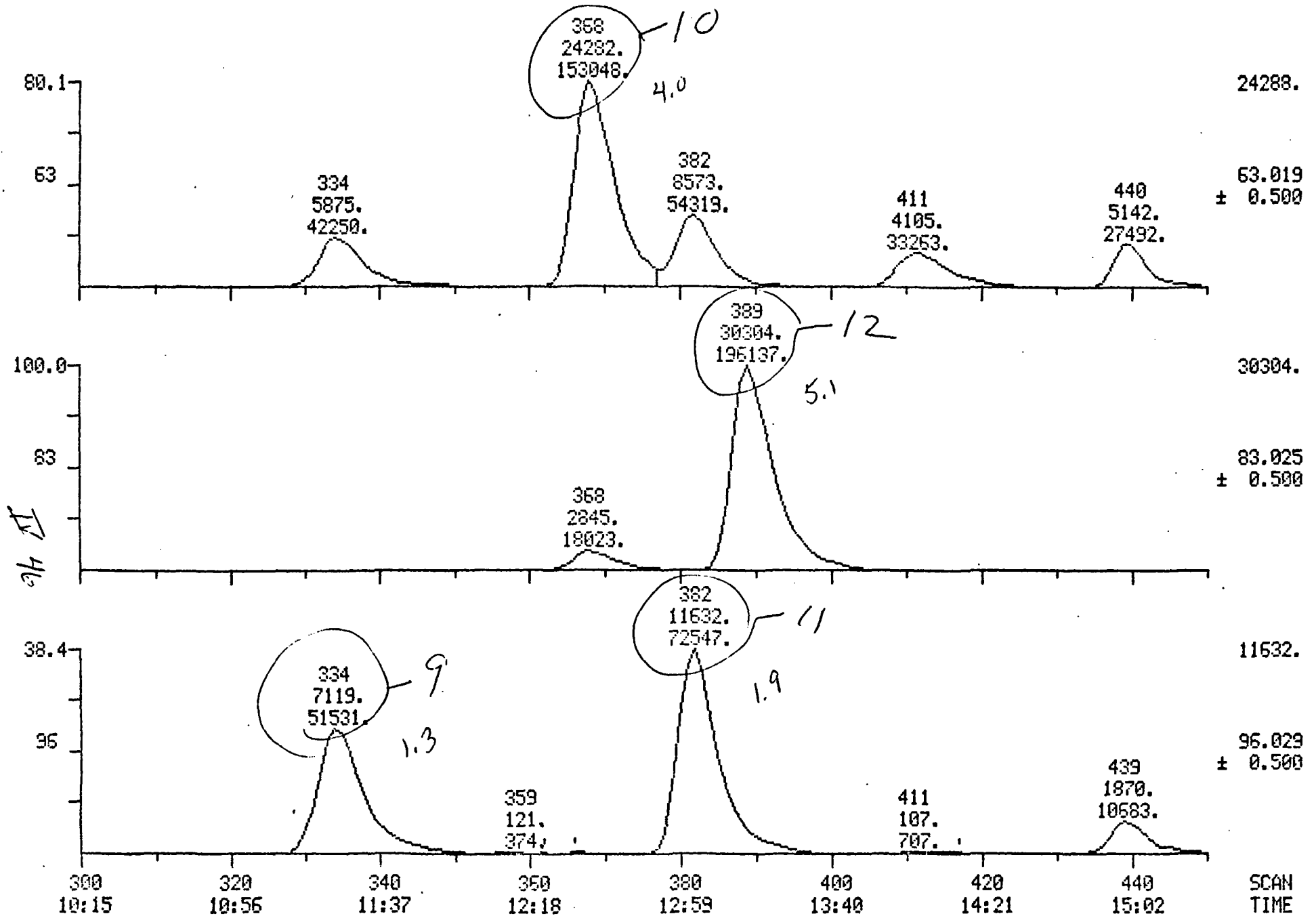
SCANS 200 TO 380



MASS CHROMATOGRAMS
03/06/85 11:43:00
SAMPLE: 50 PPB STD 3/6/85

DATA: U2755

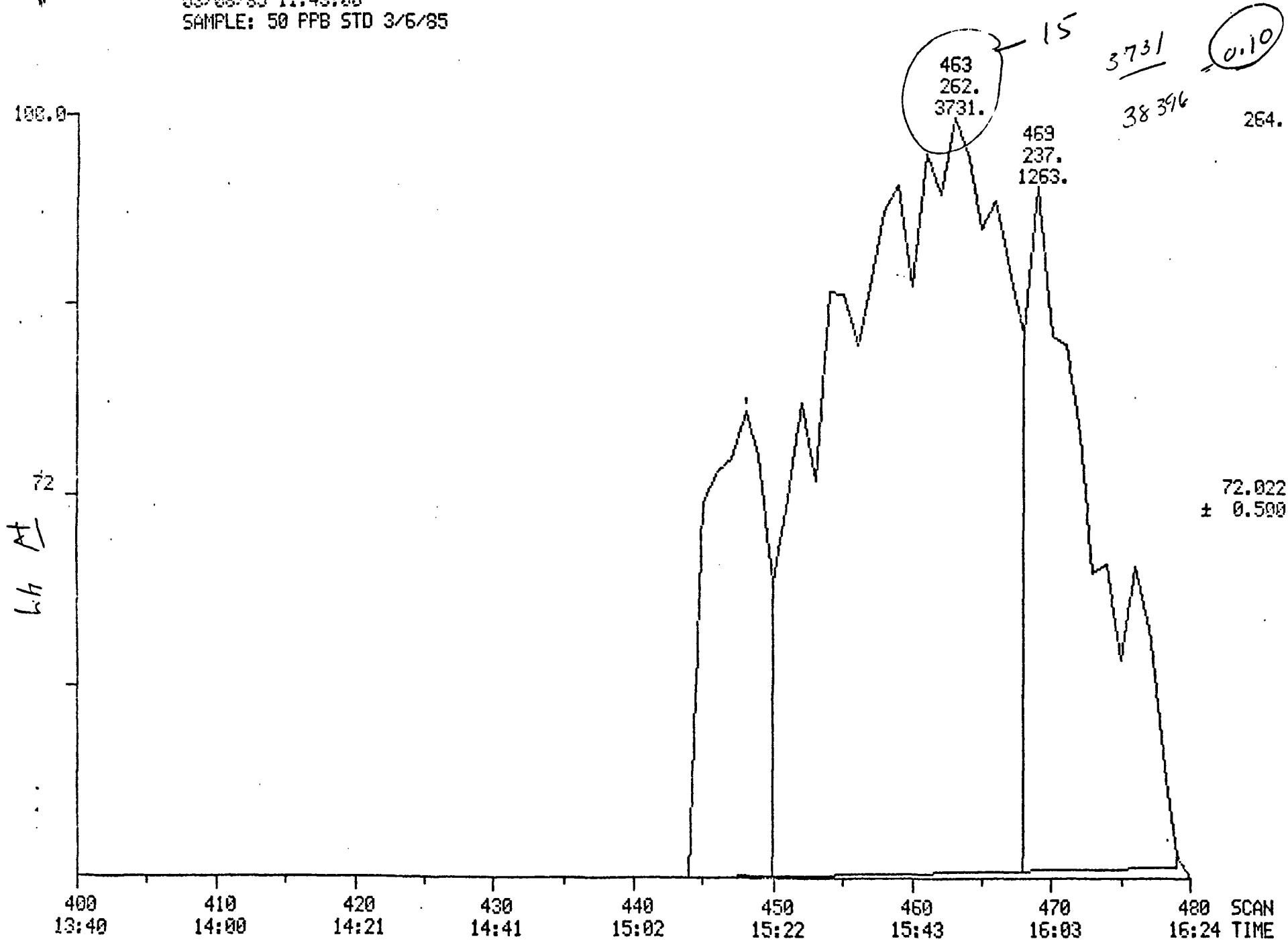
SCANS 300 TO 450



MASS CHROMATOGRAM
03/05/85 11:43:00
SAMPLE: 50 PPB STD 3/6/85

DATA: U2755

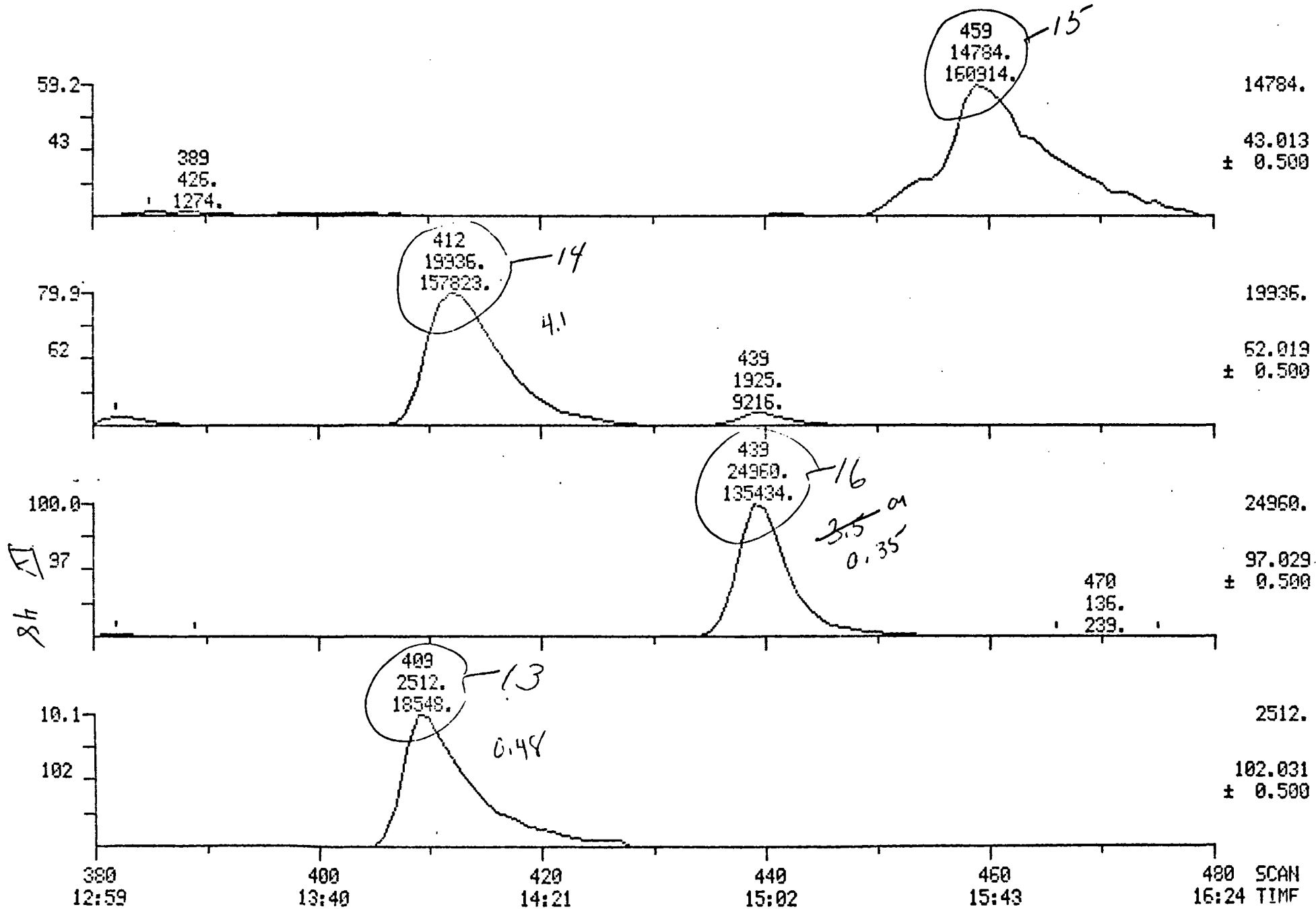
SCANS 400 TO 480



MASS CHROMATOGRAMS
03/06/85 11:43:00
SAMPLE: 50 FPB STD 3/6/85

DATA: U2755

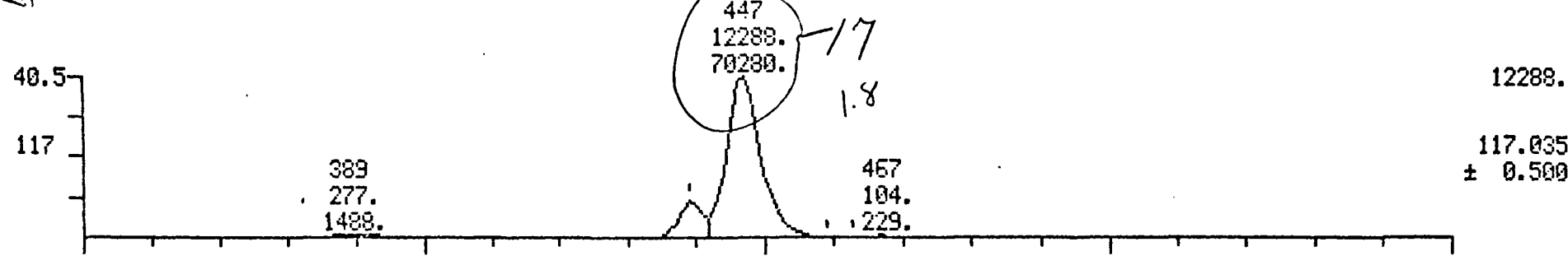
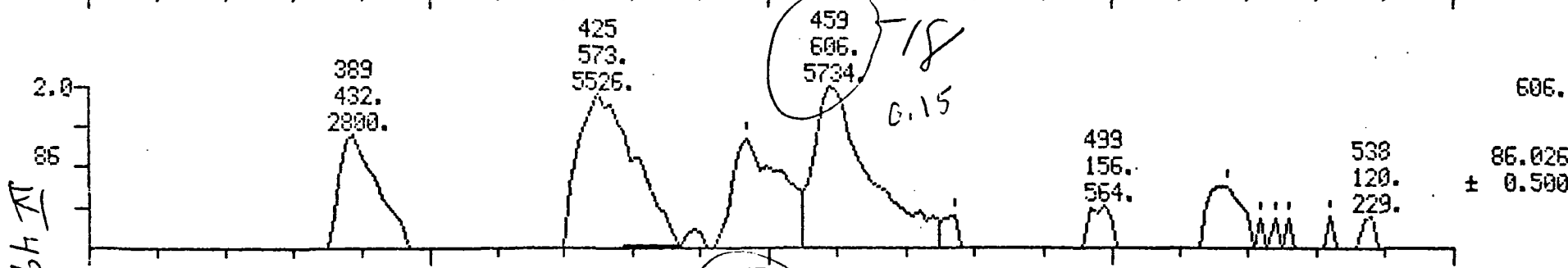
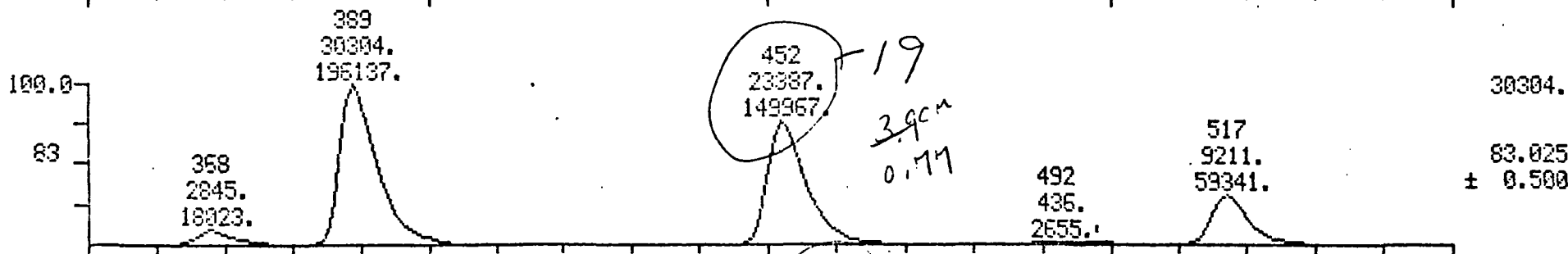
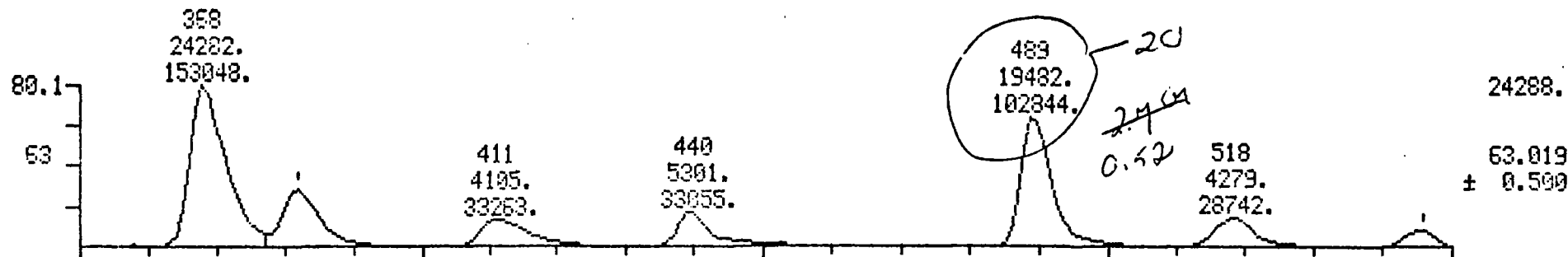
SCANS 380 TO 480



MASS CHROMATOGRAMS
 03/06/85 11:43:00
 SAMPLE: 50 PPB STD 3/6/85

DATA: U2755

SCANS 350 TO 550

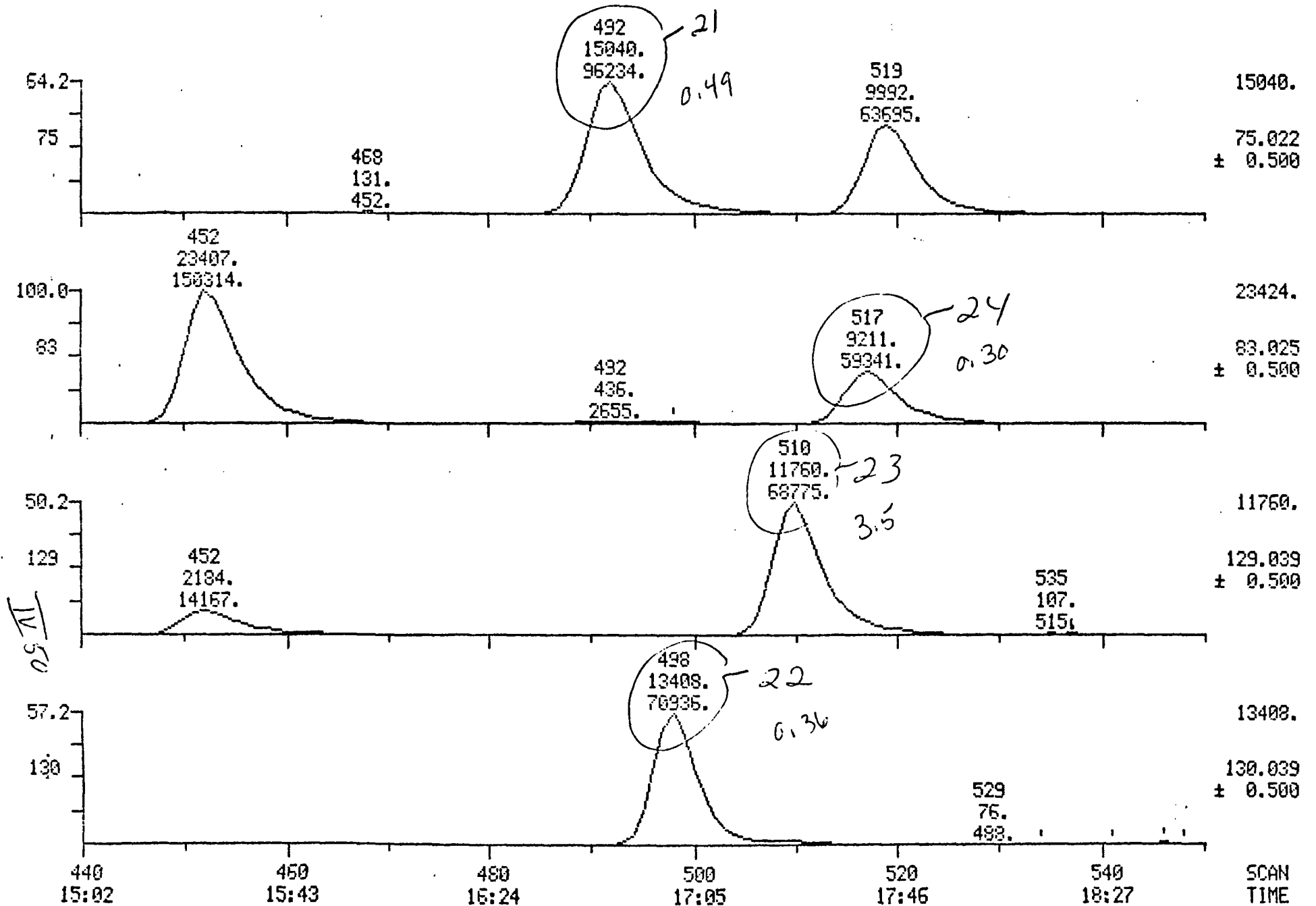


350 11:57 400 13:40 450 15:22 500 17:05 550 SCAN 18:47 TIME

MASS CHROMATOGRAMS
 03/06/85 11:43:00
 SAMPLE: 50 PFB STD 3/6/85

DATA: U2755

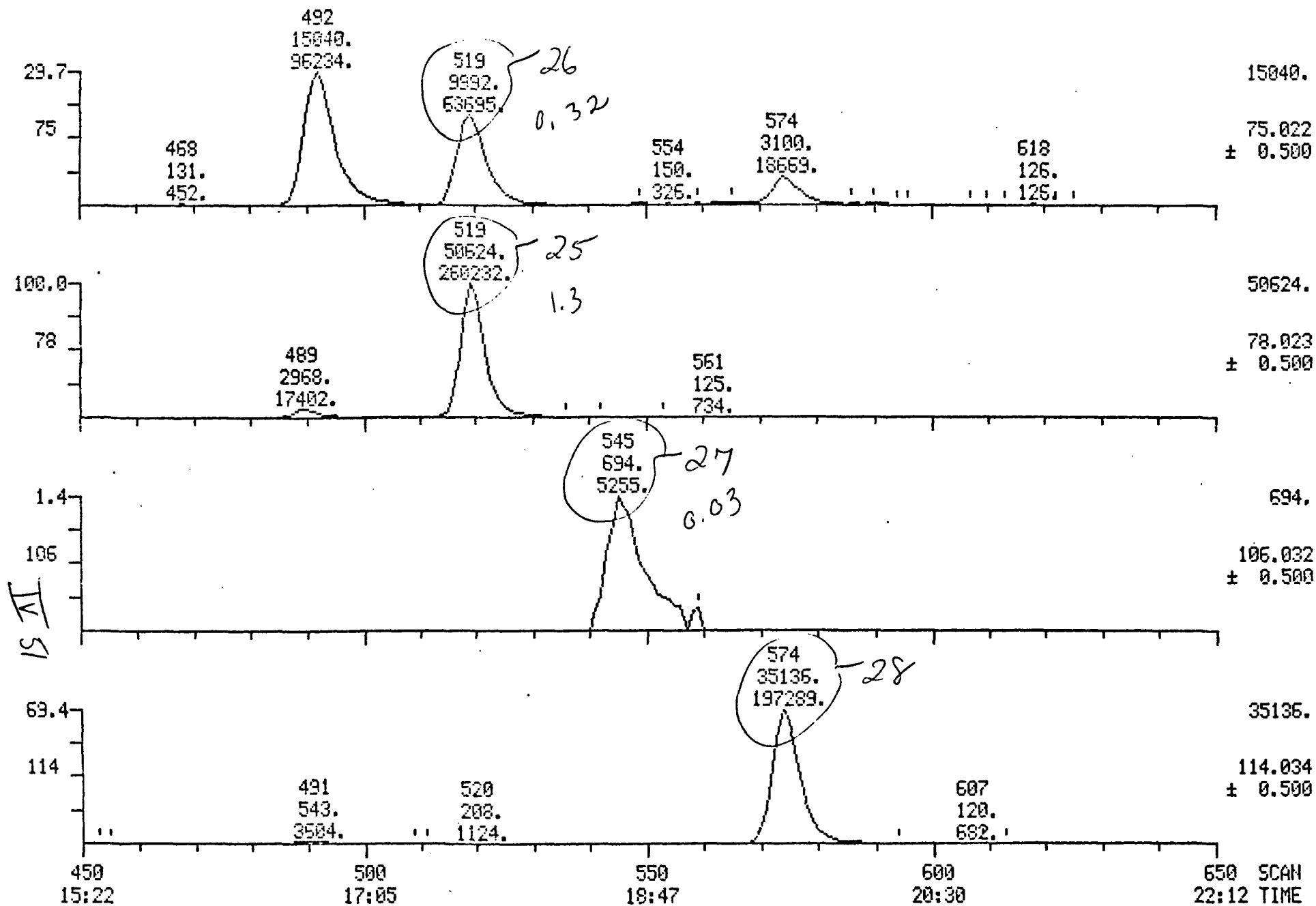
SCANS 440 TO 550



MASS CHROMATOGRAMS
03/06/85 11:43:00
SAMPLE: 50 PPB STD 3/6/85

DATA: U2755

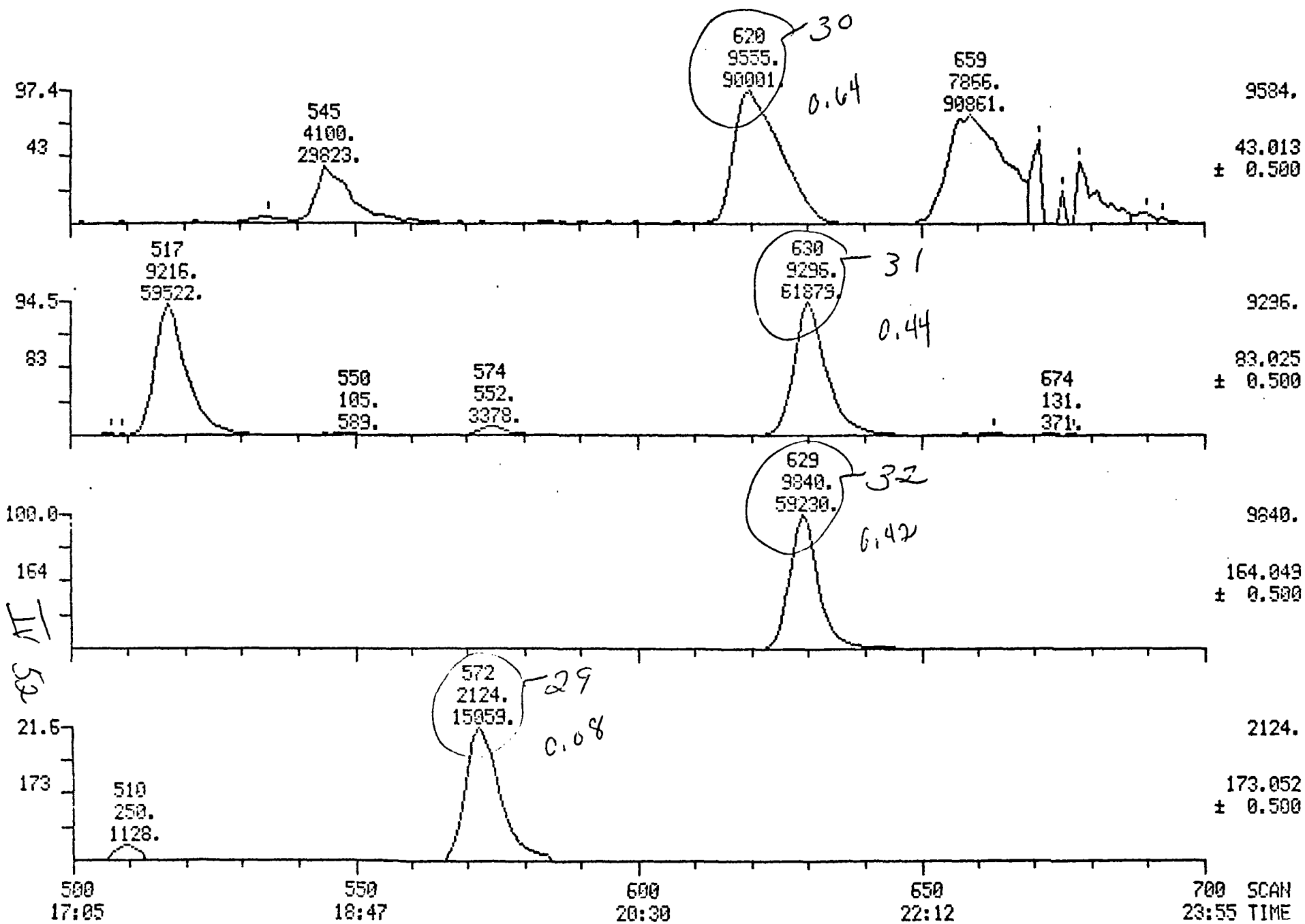
SCANS 450 TO 650



MASS CHROMATOGRAMS
03/06/85 11:43:00
SAMPLE: 50 PPB STD 3/6/85

DATA: V2755

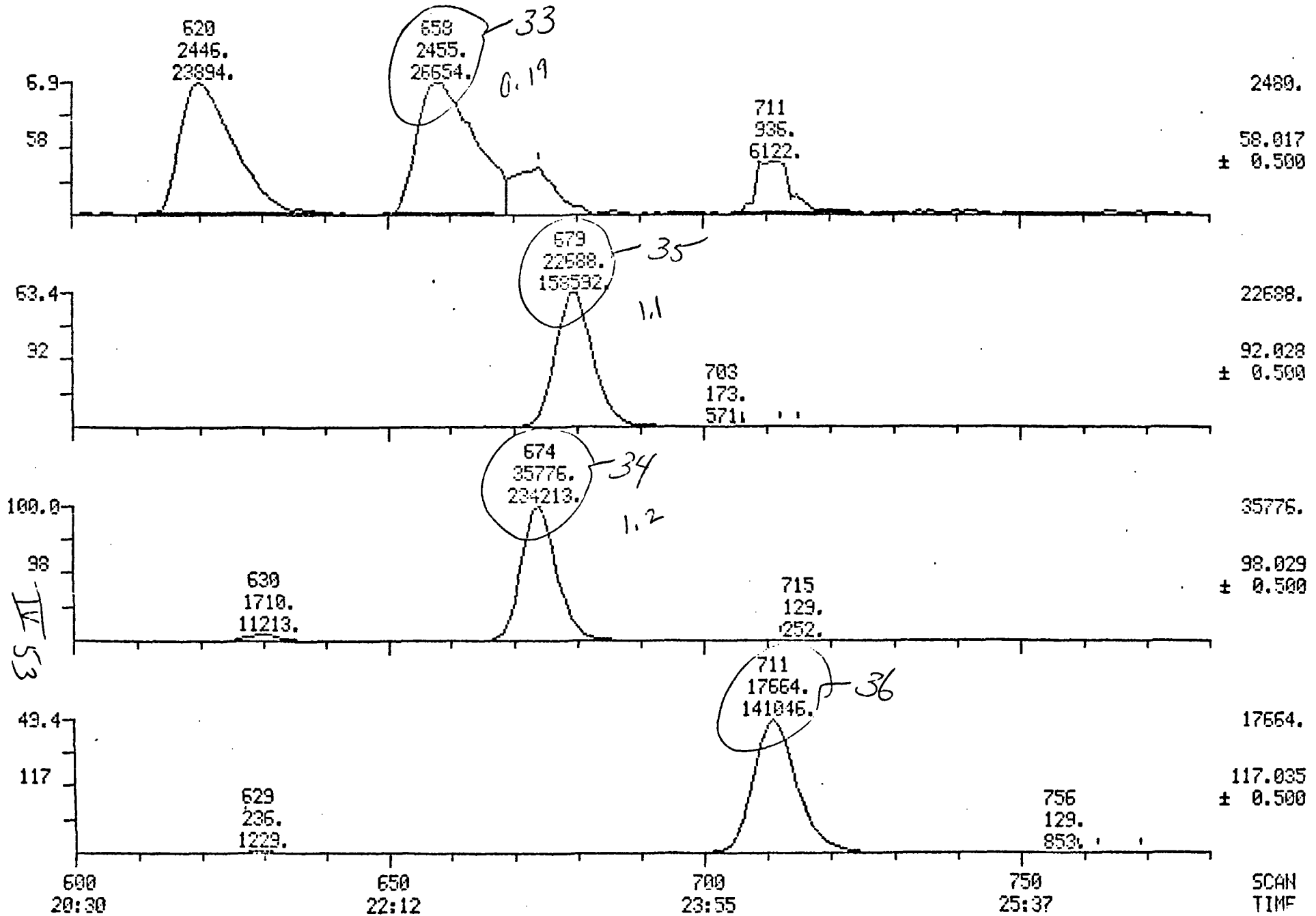
SCANS 500 TO 700



MASS CHROMATOGRAMS
 03/05/85 11:43:00
 SAMPLE: 50 PPB STD 3/6/85

DATA: U2755

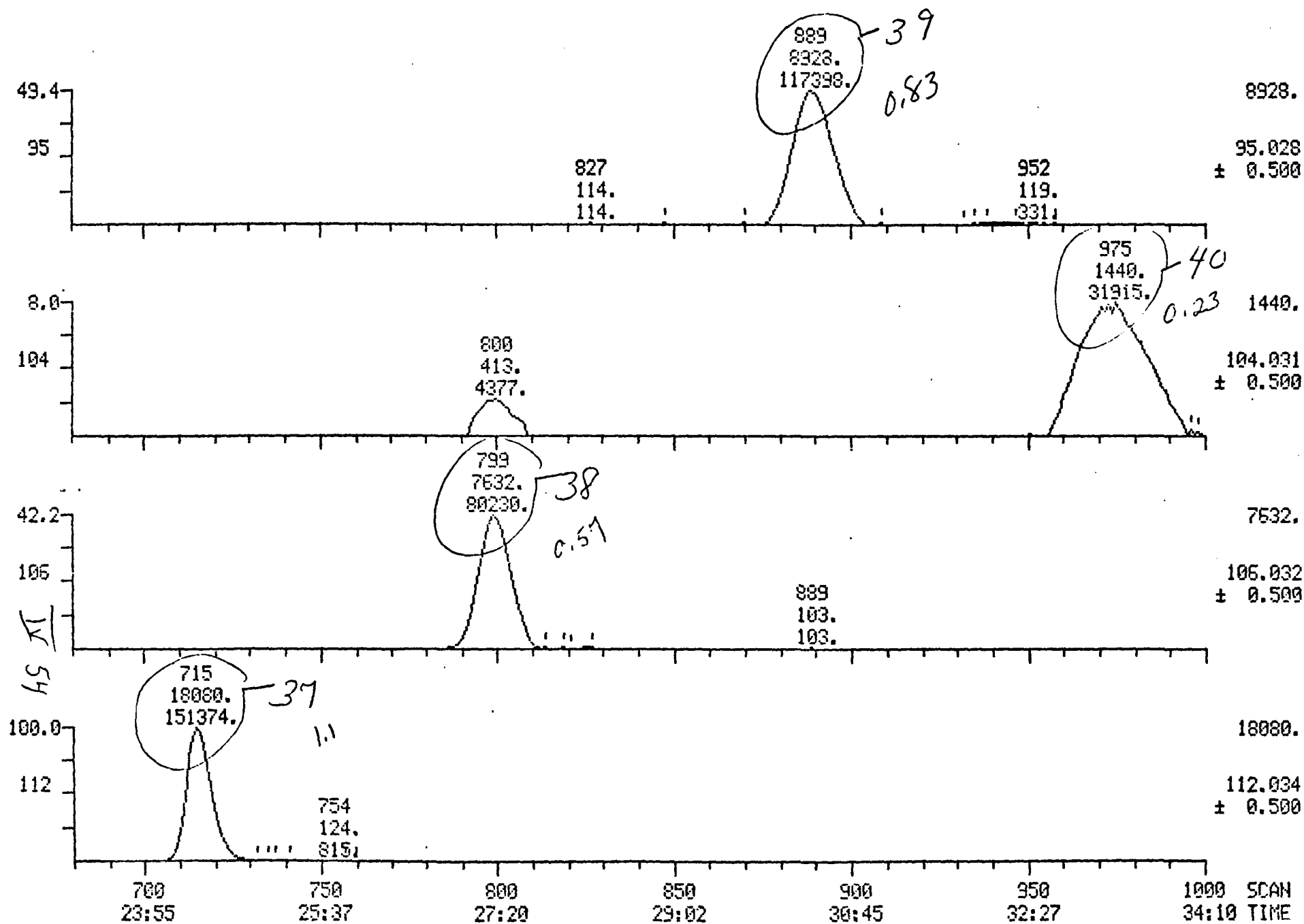
SCANS 600 TO 780



MASS CHROMATOGRAMS
03/06/85 11:43:00
SAMPLE: 50 PPB STD 3/6/85

DATA: U2755

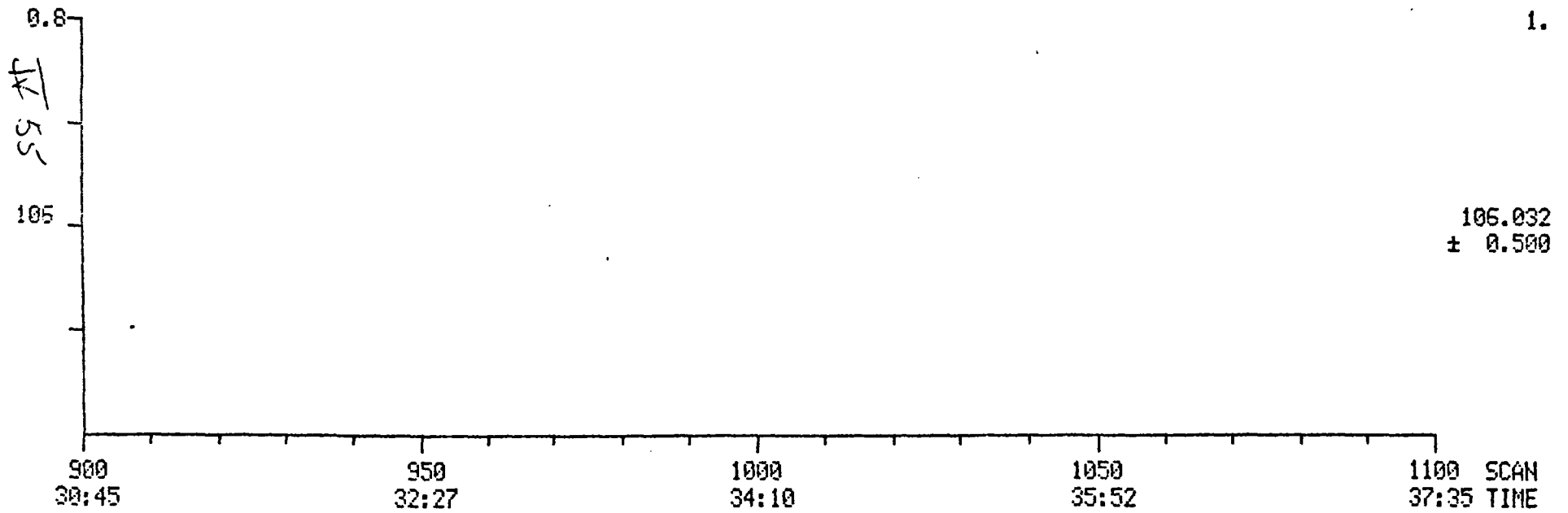
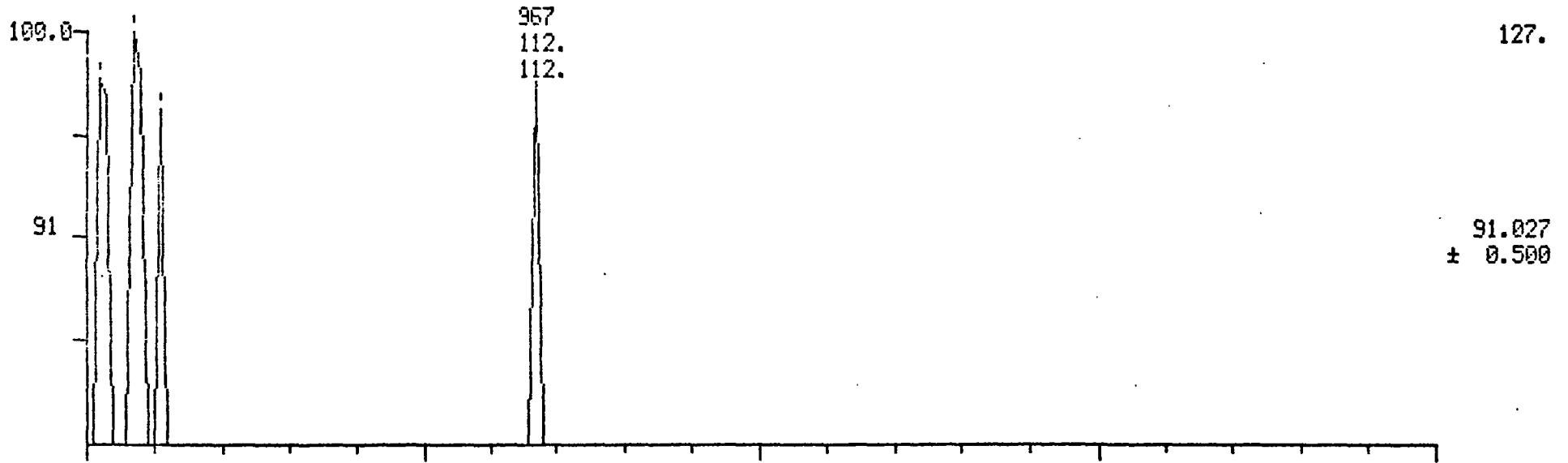
SCANS 680 TO 1000



MASS CHROMATOGRAMS
03/05/85 11:43:00
SAMPLE: 50 PPB STD 3/6/85

DATA: U2755

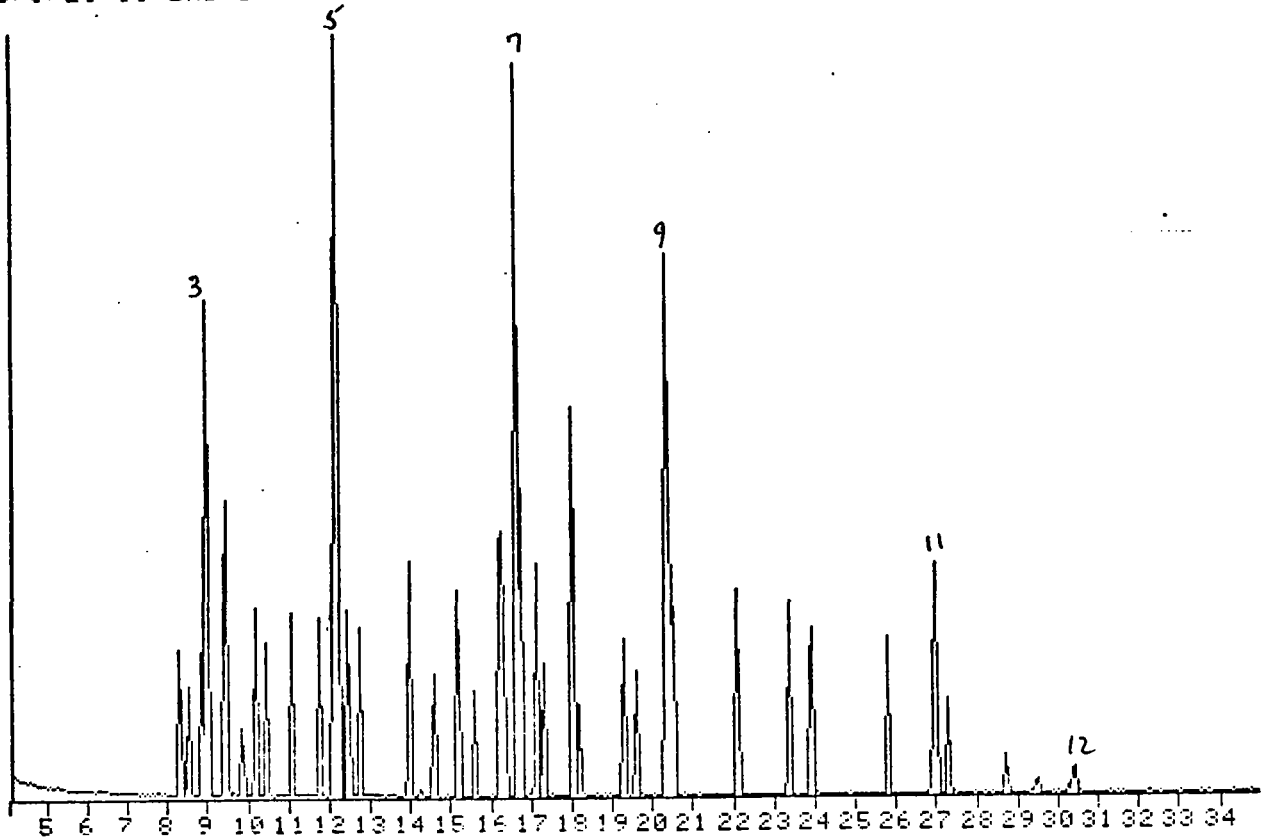
SCANS 900 TO 1100



24

74430

11e

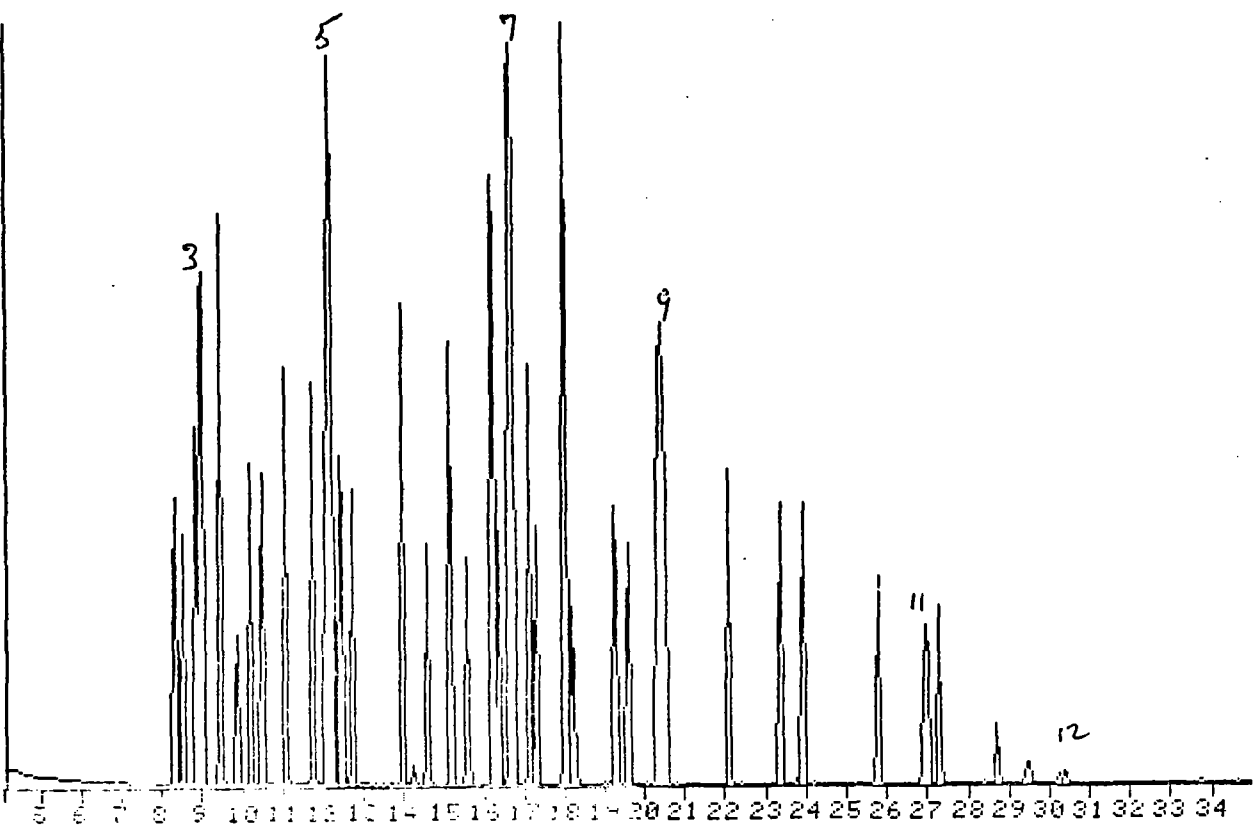


1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUORODIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV. 56

83169

ATAC

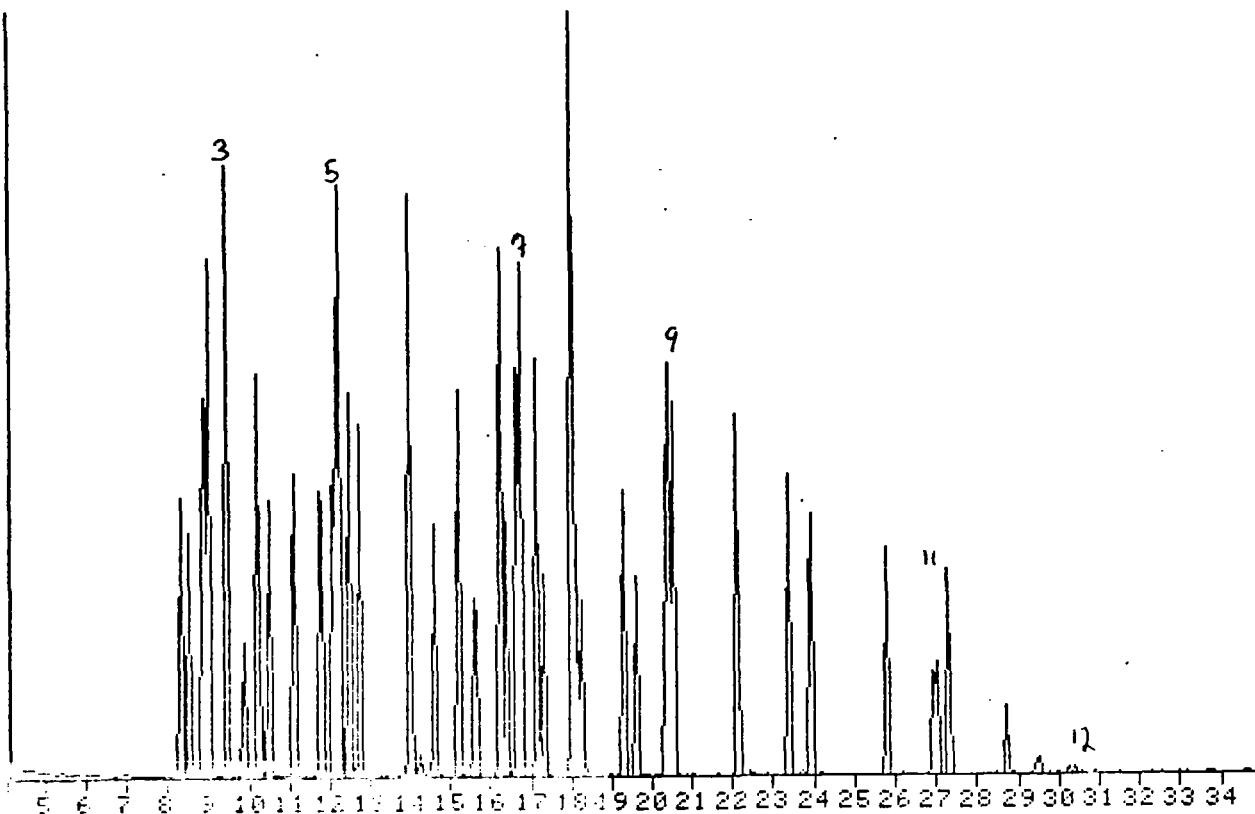


- 1. 2-FLUOROPHENOL (SURR)
- 2. PHENOL-D6 (SURR)
- 3. 1,4-DICHLOROBENZENE-D4 (IS)
- 4. NITROBENZENE-D5 (SURR)
- 5. NAPHTHALENE-D8 (IS)
- 6. 2-FLUOROBIPHENYL (SURR)
- 7. ACENAPHTHENE-I10 (IS)
- 8. 2,4,6-TRIBROMOPHENOL (SURR)
- 9. PHENANTHRENE-D10 (IS)
- 10. TERPHENYL-D14 (SURR)
- 11. CHRYSENE-D12 (IS)
- 12. PERYLENE-D12 (IS)

IV 57

132141

ATI



1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 58

80BN

3/25/85/EM

BTL#8 Q14749

D14749

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
9.0	150.0	1.000	40.0000				D4-1,4-DICHLORO BENZENE(20)
9.9	74.0	.046	.0000				N-NITROSODIMETHYLAMINE
8.4	93.0	.913	80.0000				ANILINE
8.5	93.0	.709	80.0000				BIS(2-CHLOROETHYL)ETHER
8.9	146.0	.591	80.0000				1,3-DICHLORO BENZENE
9.0	146.0	.594	80.0000				1,4-DICHLORO BENZENE
9.5	108.0	.395	80.0000				BENZYL ALCOHOL
9.5	146.0	.594	80.0000				1,2-DICHLORO BENZENE
9.9	45.0	.527	80.0000				BIS(2-CHLOROISOPROPYL)ETHER
10.2	70.0	.490	80.0000				N-NITROSODINPROPYLAMINE
10.2	117.0	.291	80.0000				HEXACHLOROETHANE
10.5	77.0	.716	80.0000				NITROBENZENE

12.2	68.0	1.000	40.0000				D8-NAPHTHALENE(20)
11.1	138.0	1.519	80.0000				ISOPHORONE
11.8	95.0	1.813	80.0000				BIS(2-CHLOROETHOXY)METHANE
12.1	180.0	2.642	80.0000				1,2,4-TRICHLORO BENZENE
12.2	129.0	1.325	80.0000				NAPHTHALENE
12.5	127.0	4.577	80.0000				4-CHLOROANILINE
12.7	225.0	1.190	80.0000				HEXACHLORO BUTADIENE
14.0	115.0	2.578	80.0000				2-METHYLNAPHTHALENE

16.7	164.0	1.000	40.0000				D10-ACENAPHTHENE
14.6	237.0	.205	80.0000				HEXACHLORO CYCLOPENTADIENE
15.2	164.0	.281	80.0000				2-CHLORONAPHTHALENE
15.6	138.0	.397	80.0000				2-NITROANILINE
16.2	164.0	.083	80.0000				DIMETHYL PHTHALATE
16.3	151.0	.272	80.0000				ACENAPHTHYLENE
16.4	165.0	.228	80.0000				2,6-DINITROTOLUENE
16.7	138.0	.293	80.0000				3-NITROANILINE
16.8	152.0	.437	80.0000				ACENAPHTHENE
17.2	139.0	.490	80.0000				DIBENZOFURAN
17.3	165.0	.320	80.0000				2,4-DINITROTOLUENE
18.0	177.0	.171	80.0000				DIETHYL PHTHALATE
18.0	165.0	.820	80.0000				FLUORENE
18.1	204.0	.358	80.0000				4-CHLOROPHENYLPHENYL ETHER
18.2	138.0	.292	80.0000				4-NITROANILINE
17.2	169.0	.144	80.0000				DIPHENYLAMINE

IV 59

3/25/85/EM

BTL#8 Q14749

D14749

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
20.3	188.0	1.000	40.0000				D10-PHENANTHRENE(20)
19.3	248.0	.112	80.0000				4-BROMOPHENYLPHENYL ETHER
19.6	284.0	.112	80.0000				HEXACHLOROBENZENE
20.4	179.0	.060	160.0000				PHENANTHRENE/ANTHRACENE
22.1	150.0	.100	80.0000				DIBUTYL PHTHALATE
23.4	202.0	.665	80.0000				FLUORANTHENE

27.0	240.0	1.000	40.0000				D12-CHRYSENE
23.9	202.0	2.264	80.0000				PYRENE
25.8	149.0	1.136	80.0000				BUTYL BENZYL PHTHALATE
26.9	228.0	.427	160.0000				BENZO(A)ANTHRACENE/CHRYSENE
27.3	149.0	1.128	80.0000				BIS(2-ETHYLHEXYL)PHTHALATE

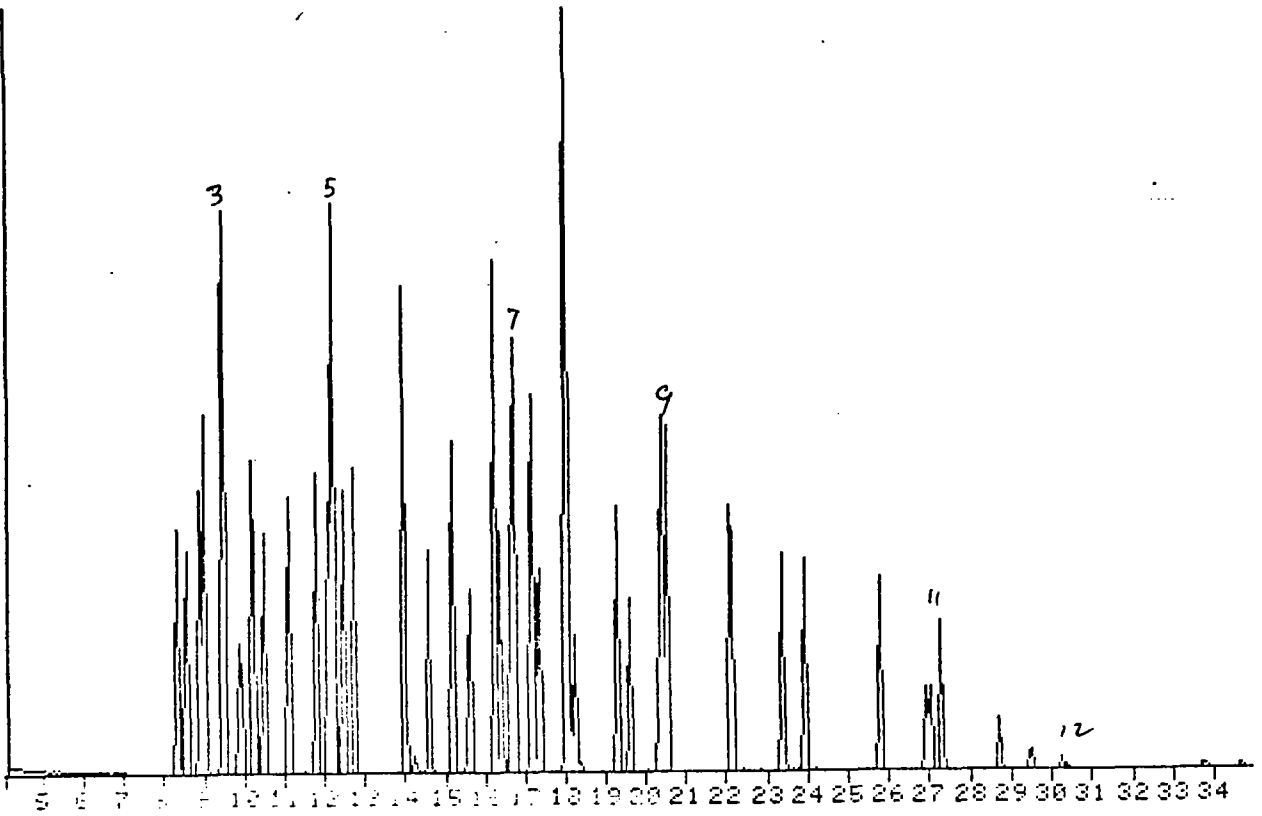
MGRE ?

30.4	264.0	1.000	40.0000				D12-PERYLENE
28.7	149.0	5.974	80.0000				DIOCTYL PHTHALATE
29.5	252.0	1.307	160.0000				BENZO(B)/(K)FLUORANTHENE
30.2	252.0	.744	80.0000				BENZO(A)PYRENE
33.7	276.0	.485	80.0000				INDENO(123-CD)PYRENE
33.8	278.0	.346	80.0000				DIBENZO(AH)ANTHRACENE
34.6	276.0	.360	80.0000				BENZO(GHI)PERYLENE

IV 60

204896

4115

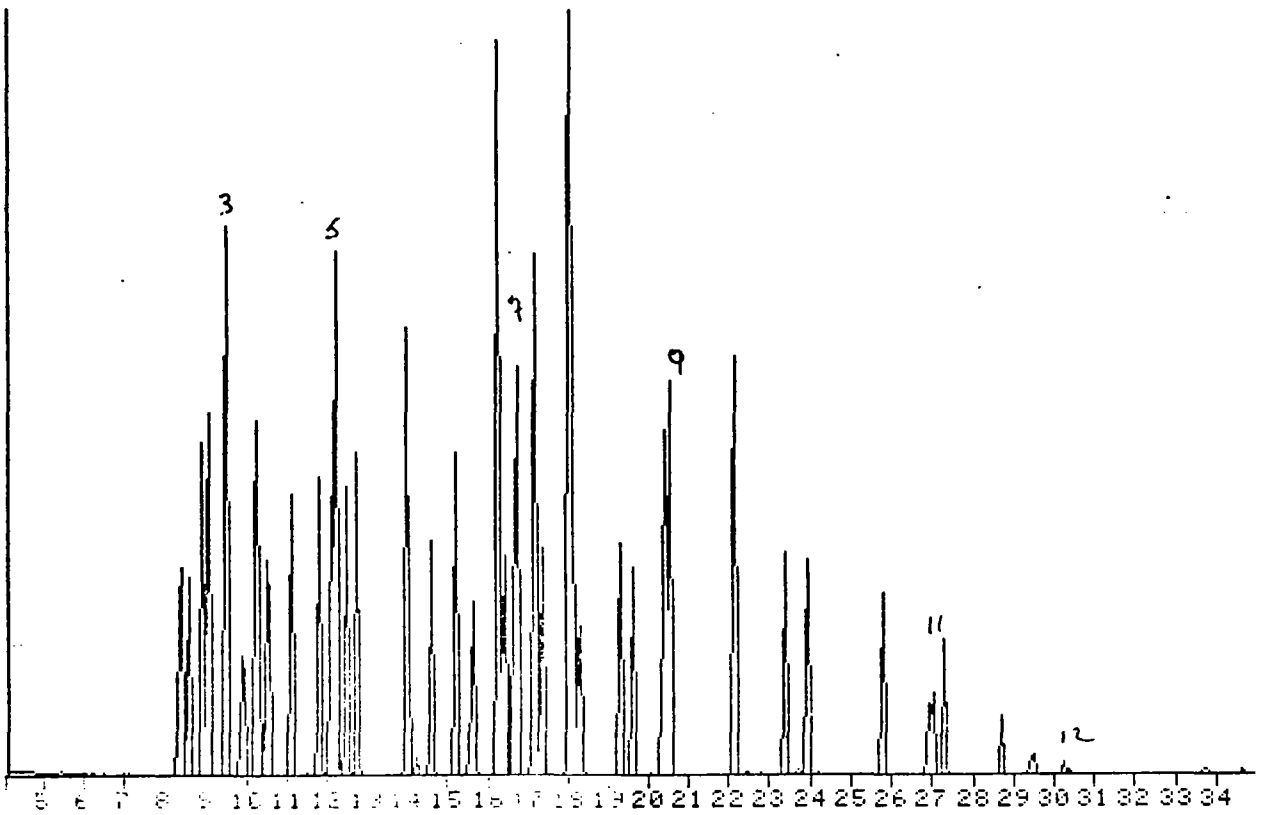


1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 61

264999

GC

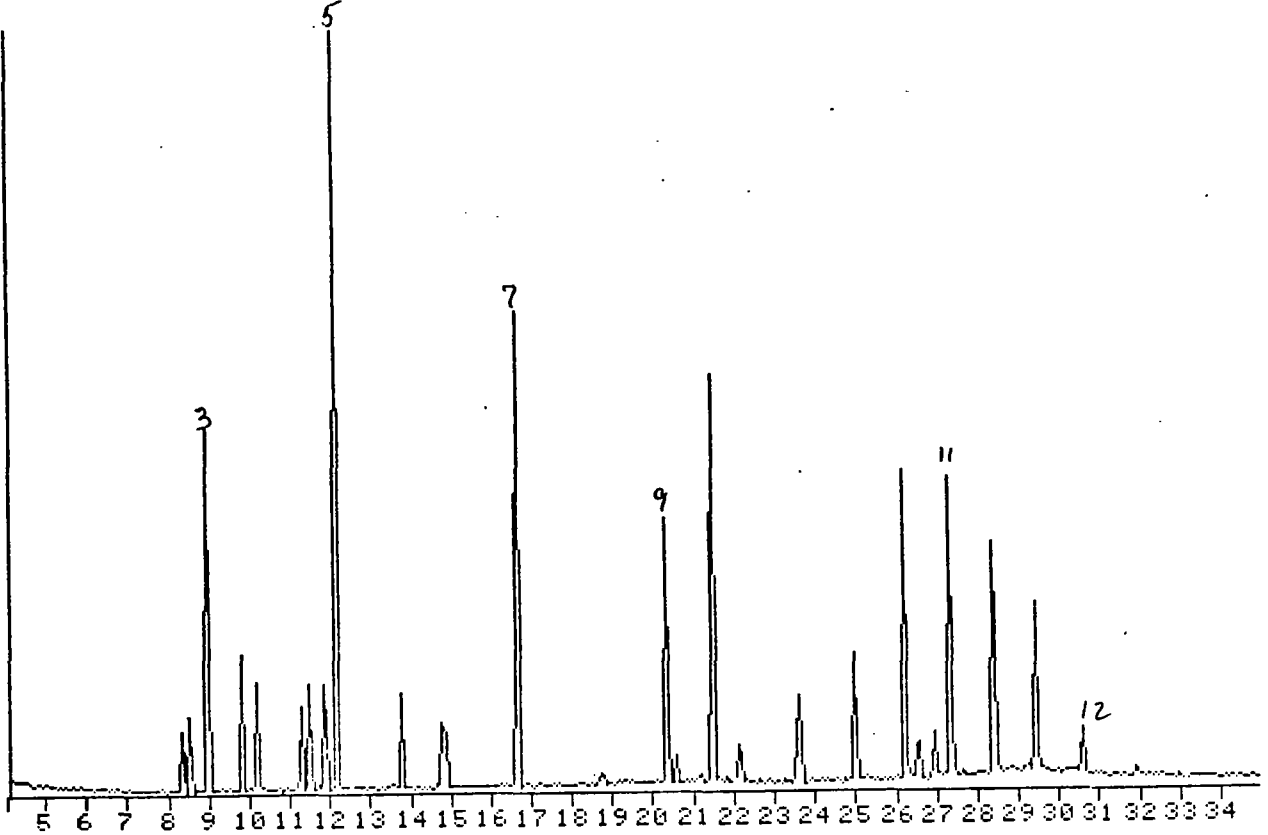


1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D16 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 62

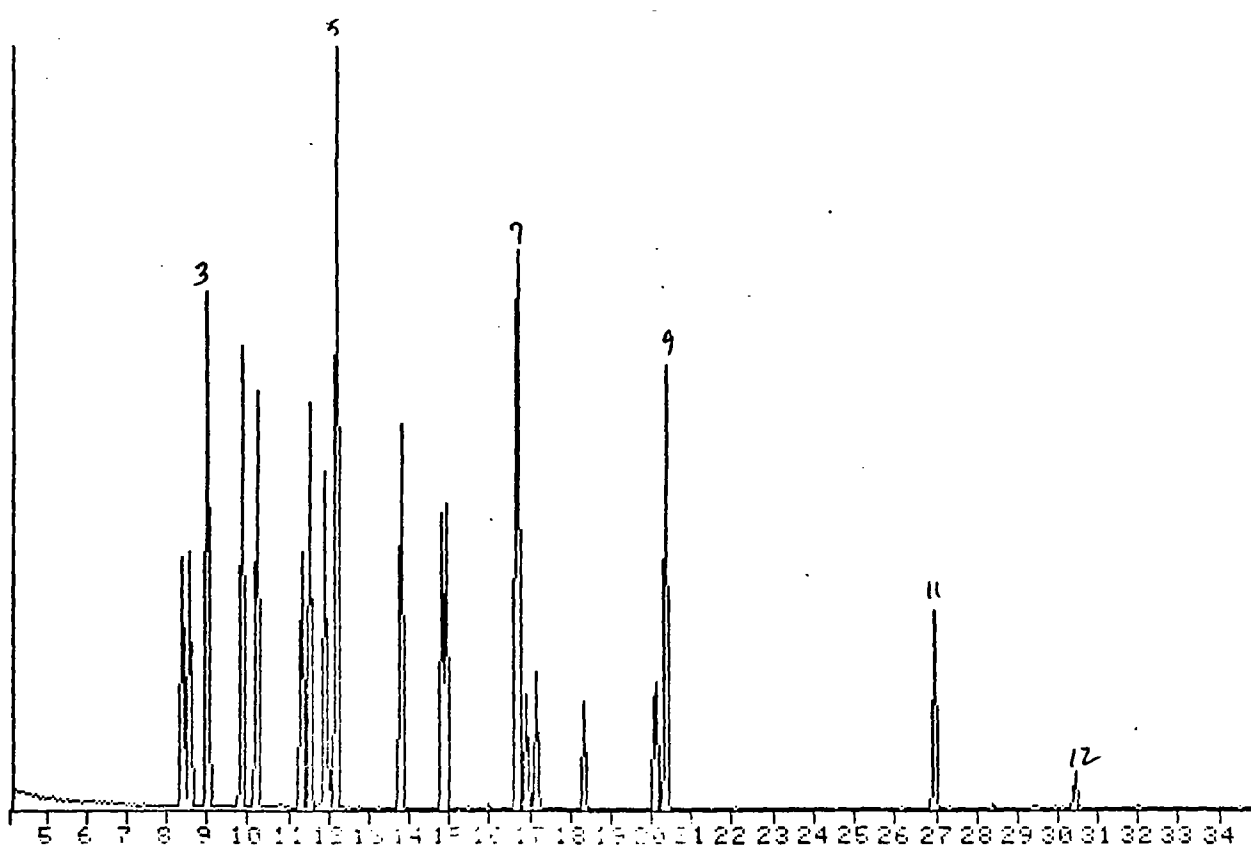
35541

4742



1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

JV 63

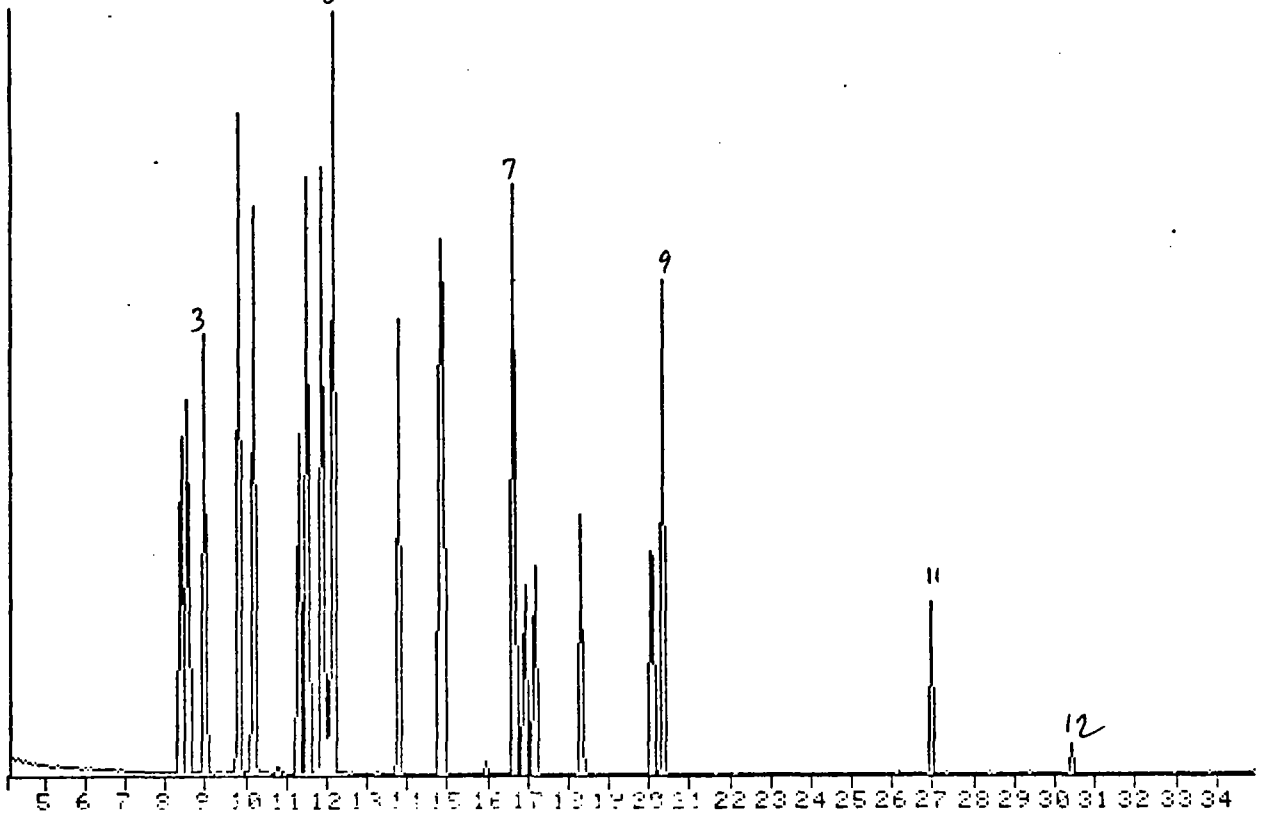


1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 64

81386

ST-10



1. 2-FLUROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUCROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 65

80A

3/25/85/EM

BTL#3 Q14744

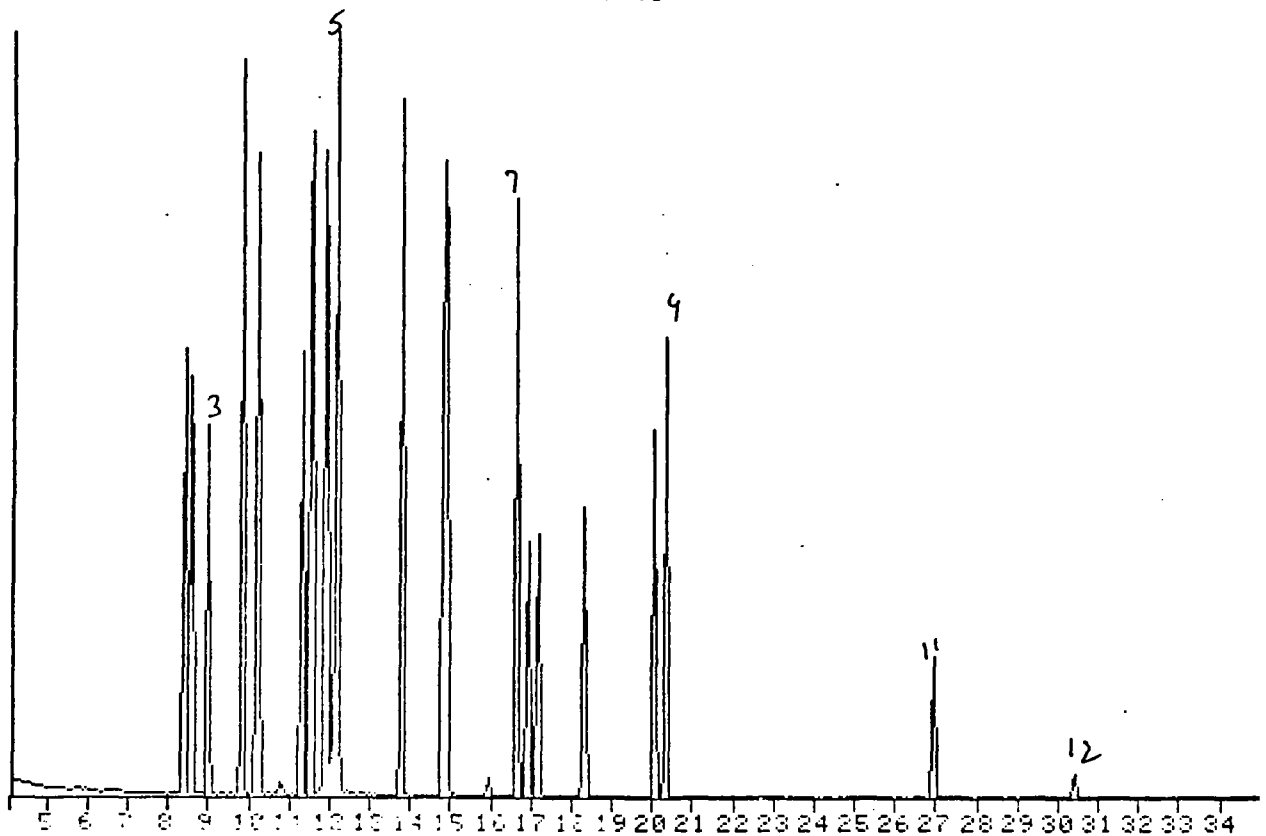
D14744

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
9.0	150.0	1.000	40.0000				D4-1,4-DICHLOROBENZENE(20)
8.4	94.0	.841	80.0000				PHENOL
8.6	128.0	.611	80.0000				2-CHLOROPHENOL
9.9	108.0	.671	80.0000				2-METHYLPHENOL
10.2	108.0	.699	80.0000				4-METHYLPHENOL
12.2	68.0	1.000	40.0000				D8-NAPHTHALENE(20)
11.3	139.0	1.759	80.0000				2-NITROPHENOL
11.5	122.0	2.581	80.0000				2,4-DIMETHYLPHENOL
12.1	122.0	1.200	80.0000				BENZOIC ACID
11.9	162.0	2.051	80.0000				2,4-DICHLOROPHENOL
13.8	142.0	2.196	80.0000				4-CHLORO-M-CRESOL
16.7	164.0	1.000	40.0000				D10-ACENAPHTHENE(20)
14.8	196.0	.237	80.0000				2,4,5-TRICHLOROPHENOL
14.8	196.0	.237	80.0000				2,4,6-TRICHLOROPHENOL
17.0	184.0	.111	80.0000				2,4-DINITROPHENOL
18.3	198.0	.126	80.0000				4,6-DINITRO-O-CRESOL
17.2	65.0	.231	80.0000				4-NITROPHENOL
20.4	188.0	1.000	40.0000				D10-PHENANTHRENE
20.1	266.0	.072	80.0000				PENTACHLOROPHENOL

IV 66

89234

318

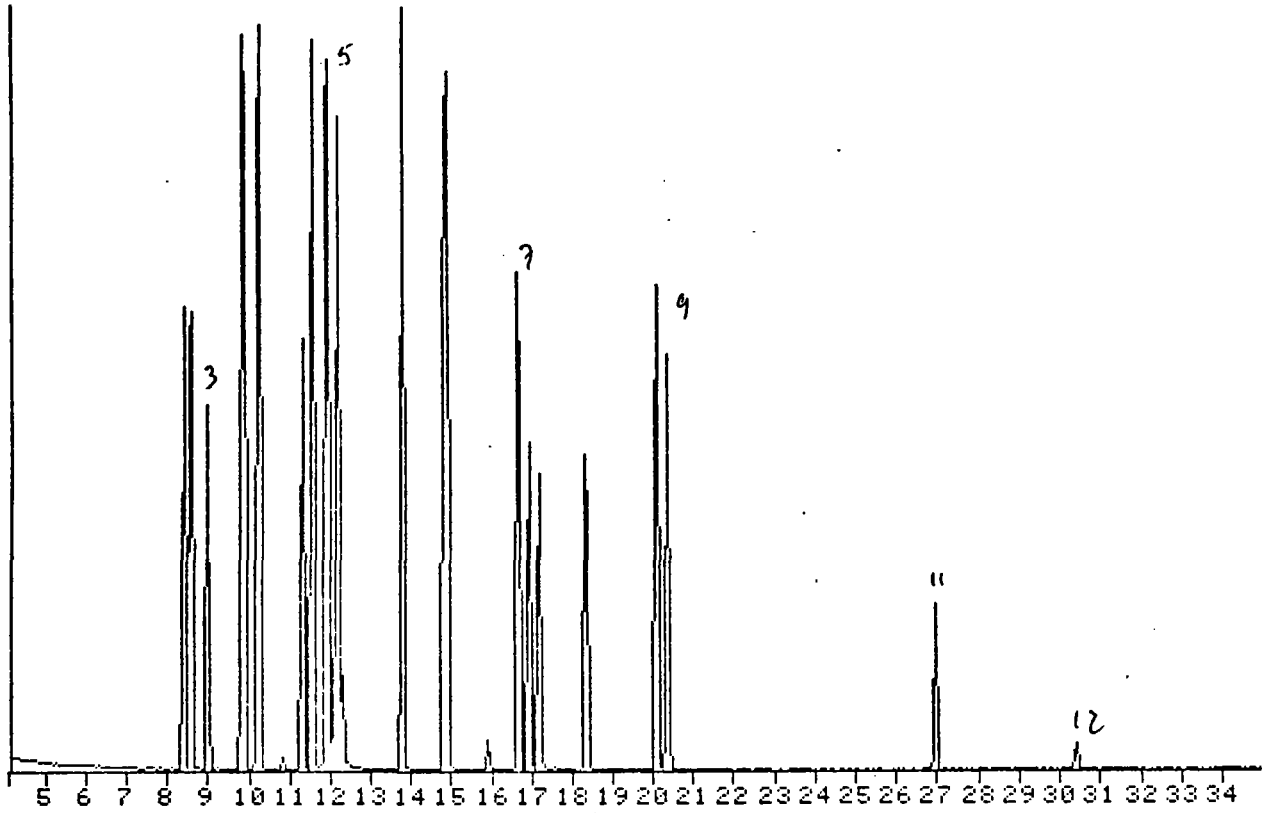


1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 67

102726

THP

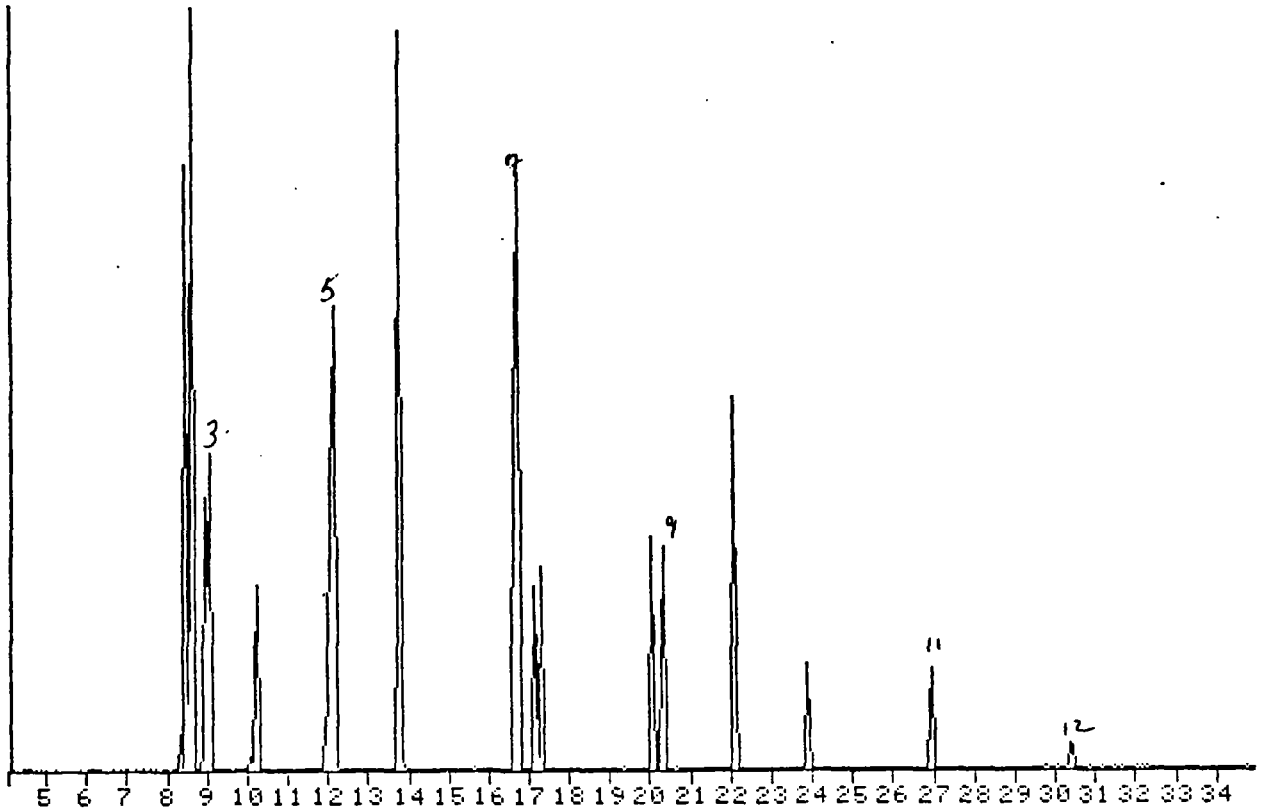


1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 68

81607

MPA



1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBIENZENE-D4 (IC)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 69

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
9.0	150.0	1.000	40.0000				D4-1,4-DICHLOROBENZENE(MS)
9.0	146.0	1.208	50.0000				1,4-DICHLOROBENZENE
10.2	70.0	1.806	50.0000				N-NITROSODINPROPYLAMINE
8.5	94.0	2.431	100.0000				PHENOL
8.6	128.0	1.819	100.0000				2-CHLOROPHENOL

12.2	136.0	1.000	40.0000				D8-NAPHTHALENE(MS)
12.1	180.0	.257	50.0000				1,2,4-TRICHLOROBENZENE
13.8	107.0	.405	100.0000				4-CHLORO-M-CRESOL

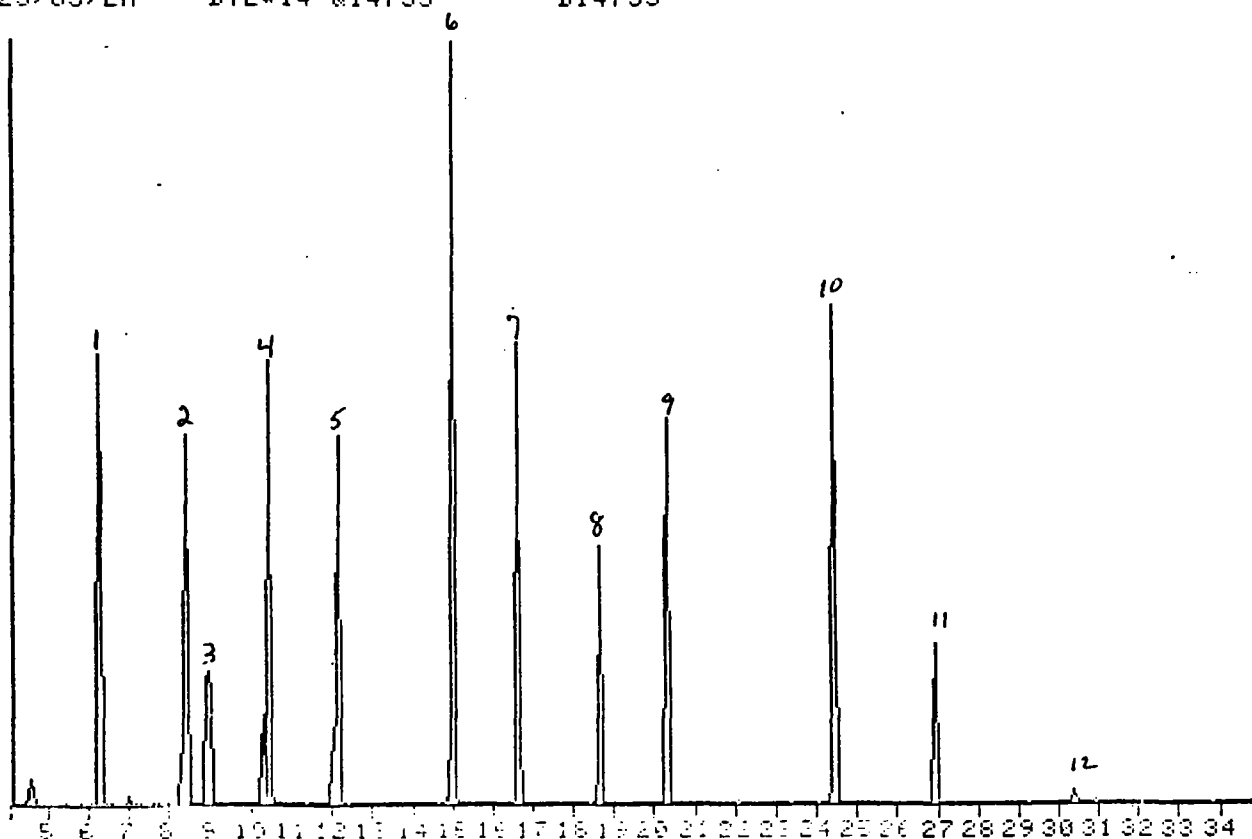
16.7	164.0	1.000	40.0000				D10-ACENAPHTHENE(MS)
16.7	153.0	1.742	50.0000				ACENAPHTHENE %R
17.3	89.0	.433	50.0000				2,4-DINITROTOLUENE %R
17.2	139.0	.251	100.0000				4-NITROPHENOL %R

20.3	188.0	1.000	40.0000				D10-PHENANTHRENE(MS)
20.1	266.0	.116	100.0000				PENTACHLOROPHENOL
22.1	149.0	2.259	50.0000				DIBUTYL PHTHALATE
23.9	202.0	.507	50.0000				PYRENE(VS ALT. IS)

IV 70

94968

700



1. 2-FLUOROPHENOL (SUPP)
2. PHENOL-D6 (SUPP)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SUPP)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROPHENYL (SUPP)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SUPP)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SUPP)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

10 71

90

SS-9

3/25/85/EM

BTL#14 Q14755

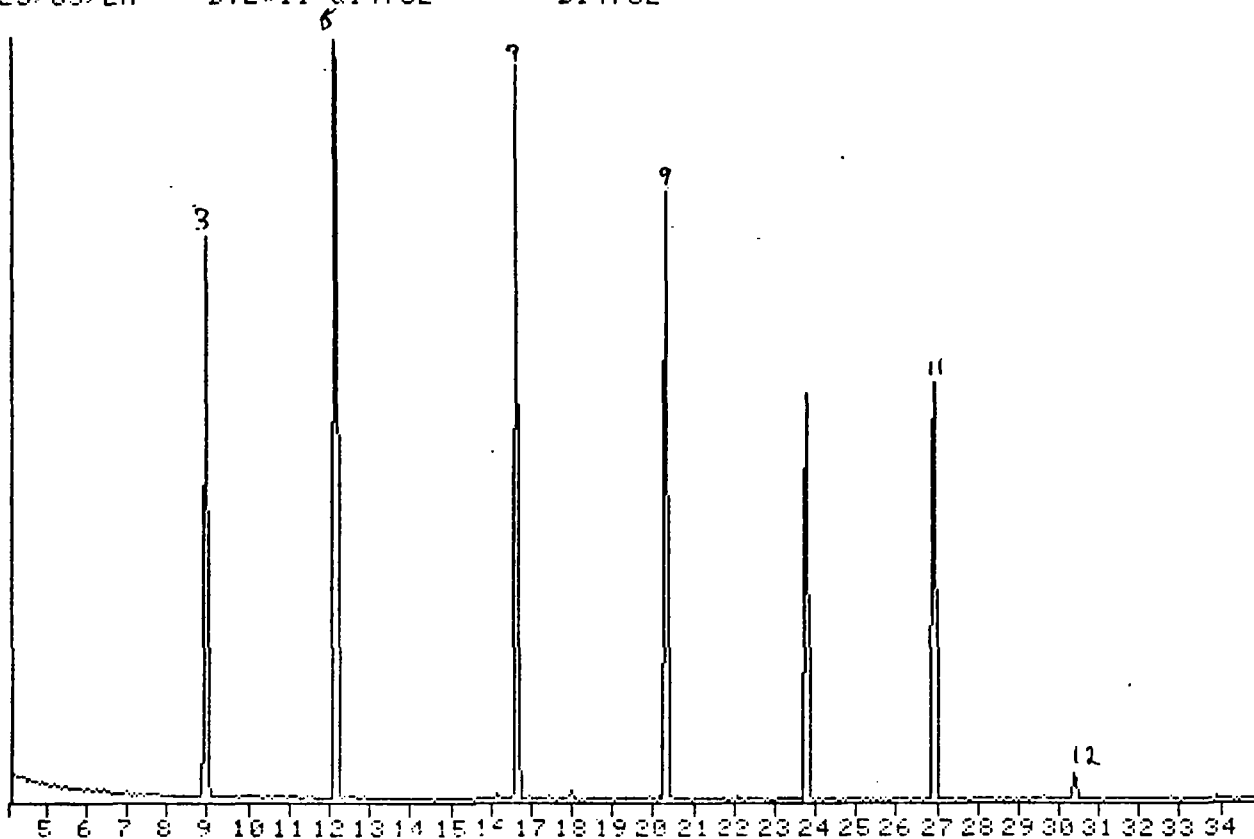
D14755

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
8.9	150.0	1.000	40.0000				D4-1,4-DICHLOROBENZENE
8.4	99.0	1.343	100.0000				D6-PHENOL
6.3	112.0	1.043	100.0000				2-FLUROPHENOL%R
16.7	164.0	1.000	40.0000				D10-ACENAPHTHENE
10.5	128.0	.310	100.0000				D5 NITROBENZENE
16.7	164.0	1.000	40.0000				D10-ACENAPHTHENE
15.0	172.0	.996	100.0000				2-FLUOROBIPHENYL
20.3	188.0	1.000	40.0000				D10-PHENANTHRENE
18.7	330.0	.053	100.0000				TRIBROMOPHENOL %R
27.0	240.0	1.000	40.0000				D-12 CHRYSENE
24.5	244.0	1.182	100.0000				D14-TERPHENYL %R

15 72

59128

STIC



1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 73

FILE NUMBER 14752

100 B/DCLB

3/25/85/EM

BTL#11 Q14752

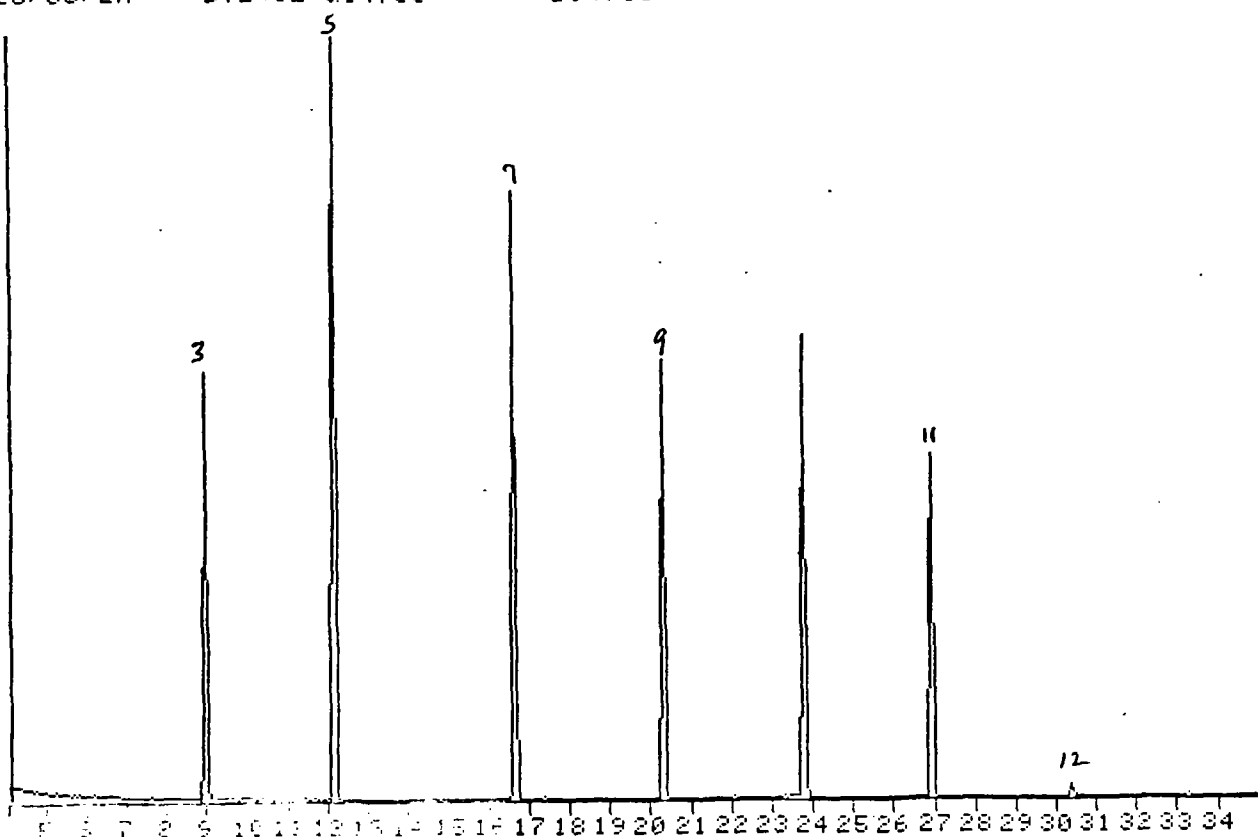
D14752

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
27.0	240.0	1.000	40.0000				D12-CHRYSENE
23.8	184.0	.712	100.0000				BENZIDINE
27.0	252.0	.146	100.0000				DICHLOROBENZIDINE

IV 74

85220

14753



1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUORODIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

IV 75

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
27.0	240.0	1.000	40.0000				D12-CHRYSENE
23.8	184.0	.839	200.0000				BENZIDINE
27.0	252.0	.099	200.0000				DICHLOROBENZIDINE

IV 76

OK

Case No. 3967 Laboratory GC A
 Contract No. 5705-050 GC Column SP 2250/240 GC Instrument ID MP 5780
6801-6767

DATE OF ANALYSIS 3/30/85 DATE OF ANALYSIS 3/30/85
 TIME OF ANALYSIS 12:52, 13:48 TIME OF ANALYSIS 3:36, 4:31
 LABORATORY ID P-30-4, P-31-4 LABORATORY ID P-30-3, P-31-3

COMPOUND	RT	RETENTION TIME WINDOW	CALIBRATION FACTOR	CONF. OR QUANT.	RT	CALIBRATION FACTOR	CONF. OR QUANT.	PERCENT DIFF. **
alpha-BHC	2.01	1.99-2.01	12.57/05 = 312	Conf.	2.00	58.64/125 = 469	Conf.	40. a
beta-BHC	3.03	3.01-3.03	10.77/05 = 215	Conf.	3.02	27.22/125 = 218	Conf.	1.8
delta-BHC	3.54	3.52-3.56	13.54/05 = 272	Conf.	3.54	40.74/125 = 326	Conf.	17.
gamma-BHC	2.59	2.56-2.60	14.10/05 = 282	Conf.	2.58	57.54/125 = 460	Conf.	35. a
Heptachlor	3.19	3.16-3.20	11.12/05 = 222	Conf.	3.18	56.57/125 = 453	Conf.	18.
Aldrin	3.84	3.82-3.89	14.74/05 = 295	Conf.	3.88	46.82/125 = 374	Conf.	23. a
Heptachlor Epoxide	6.03	6.01-6.03	18.44/05 = 369	Conf.	6.02	41.01/125 = 328	Conf.	6.6
Endosulfan I	7.65	7.61-7.67	13.07/05 = 261	Conf.	7.64	45.52/125 = 364	Conf.	2.4
Dieldrin	9.47	9.41-9.49	22.57/05 = 451	Conf.	9.44	36.47/125 = 292	Conf.	16.
4,4'-DDE	9.14	9.12-9.14	16.17/05 = 323	Conf.	9.13	42.47/125 = 339	Conf.	2.9
Endrin	11.57	11.51-11.59	26.47/05 = 529	Conf.	11.54	26.27/125 = 210	Conf.	4.6
Endosulfan II	14.22	14.14-14.24	16.17/05 = 323	Conf.	14.19	40.44/125 = 324	Conf.	0.31
4,4'-DDD	14.39	14.35-14.41	7.73/05 = 155	Conf.	14.36	21.57/125 = 172	Conf.	10.
Endrin Aldehyde	18.74	18.77-18.83	13.21/05 = 264	Conf.	18.74	41.26/125 = 330	Conf.	18.
Endosulfan Sulfate	23.30	23.11-23.35	13.53/05 = 271	Conf.	23.23	29.63/125 = 237	Conf.	12.
4,4'-DDT	17.42	17.33-17.45	5.44/05 = 109	Conf.	17.37	16.14/125 = 129	Conf.	18.
Methoxychlor	34.80*	34.76-34.88	7.07/125 = 56.56*	Conf.	34.84*	2.14/25 = 8.56*	Conf.	17.
Endrin Ketone	31.73	31.55-31.71	14.47/05 = 289	Conf.	31.65	25.50/125 = 204	Conf.	8.2
Techn. Chlordane								
alpha-Chlordane								
gamma-Chlordane								
Toxaphene								
Arceclor - 1016								
Arceclor - 1221								
Arceclor - 1232								
Arceclor - 1242								
Arceclor - 1248								
Arceclor - 1254								
Arceclor - 1260								

a See accompanying sheet for comparison of duplicate injections of the same standard which resulted in good agreement of calibration factors. This indicates that there was a problem with the standard solutions and not an instrument problem.

IV
79

* SEE EXHIBIT B PART 7 * from P-30-3
 a. Does peak integrate due to crossover B.L.

** CONF. = CONFIRMATION (LOW DIFFERENCE) MEAN CALIBRATION (LOW DIFFERENCE)

Case No. 3964
 Contract No. 3-765-050 PN
68-01-6767

Laboratory GCA
 GC Column SP 2250/2401 GC Instrument ID HP 5770

DATE OF ANALYSIS _____	DATE OF ANALYSIS <u>3/30</u>
TIME OF ANALYSIS _____	TIME OF ANALYSIS <u>10:57</u>
LABORATORY ID _____	LABORATORY ID <u>P-30-4</u>

COMPOUND	RT	RETENTION TIME WINDOW	CALIBRATION FACTOR	CONF. OR QUANT.	RT	CALIBRATION FACTOR	CONF. OR QUANT.	PERCENT DIFF.**
alpha-BHC					2.00	15.65/0.5 = 312		0
beta-BHC								
delta-BHC								
gamma-BHC					2.58	14.74/0.5 = 294		1.0
Heptachlor								
Aldrin					3.87	15.07/0.5 = 301		1.7
Heptachlor Epoxide								
Endosulfan I								
Dieldrin								
4,4'-DDE								
Endrin								
Endosulfan II								
4,4'-DDD								
Endrin Aldehyde								
Endosulfan Sulfate								
4,4'-DDT								
Methoxychlor								
Endrin Ketone								
Tech. Chlordane								
alpha-Chlordane ³								
gamma-Chlordane ⁴								
Texaphene								
Aroclor - 1016								
Aroclor - 1221								
Aroclor - 1232								
Aroclor - 1242								
Aroclor - 1248								
Aroclor - 1254								
Aroclor - 1260								

IV
878

* SEE EXHIBIT B, PART 7

** CONF. = CONFIRMATION (<20% DIFFERENCE)
 QUANT. = QUANTITATION (<10% DIFFERENCE)

Case No. 5169 Laboratory 6CA
 Contract No. 3-705-050 PN GC Column SP 2250/2401 GC Instrument ID HP5880
68-01-6767

DATE OF ANALYSIS 4/1/85
 TIME OF ANALYSIS 14:23, 15:05
 LABORATORY ID P-30-4, P 31-4

DATE OF ANALYSIS 4/2/85
 TIME OF ANALYSIS 4:16, 4:57
 LABORATORY ID P-30-4, P 31-4

COMPOUND	RT	RETENTION TIME WINDOW	CALIBRATION FACTOR	CONF. OR QUANT.	RT	CALIBRATION FACTOR	CONF. OR QUANT.	PERCENT DIFF. **
alpha-BHC	2.00	1.79-2.01	24.11/05 = 476	Conf.	2.00	24.71/05 = 496	Conf.	4.1
beta-BHC	3.02	2.99-3.05	17.00/05 = 174	Conf.	3.02	15.61/05 = 170	Conf.	3.2
delta-BHC	3.53	3.50-3.58	16.71/05 = 216	Conf.	3.53	11.61/05 = 234	Conf.	8.0
gamma-BHC	2.58	2.57-2.59	21.75/05 = 435	Conf.	2.58	22.67/05 = 452	Conf.	3.8
Heptachlor	3.18	3.17-3.19	28.11/05 = 563	Conf.	3.18	28.11/05 = 562	Conf.	0.18
Aldrin	3.88	3.86-3.90	21.14/05 = 423	Conf.	3.87	21.77/05 = 435	Conf.	2.5
Heptachlor Epoxide	6.02	5.97-6.05	24.16/05 = 482	Conf.	6.01	24.51/05 = 487	Conf.	1.0
Endosulfan I	7.63	7.56-7.70	16.03/05 = 321	Conf.	7.62	16.01/05 = 324	Conf.	0.73
Dieldrin	9.43	9.41-9.45	17.46/05 = 349	Conf.	9.42	20.27/05 = 405	Conf.	15.0
4,4'-DDE	9.12	9.04-9.22	13.90/05 = 278	Conf.	9.11	14.33/05 = 287	Conf.	3.2
Endrin	11.53	11.49-11.55	17.00/05 = 249	Conf.	11.51	14.71/05 = 294	Conf.	17.0
Endosulfan II	14.17	14.07-14.28	14.26/05 = 285	Conf.	14.15	14.29/05 = 286	Conf.	0.35
4,4'-DDD	14.35	14.28-14.40	10.54/05 = 206	Conf.	14.33	14.64/05 = 293	Conf.	35.0
Endrin Acidhyde	18.89	18.83-18.95	20.12/05 = 402	Conf.	18.85	15.73/05 = 314	Conf.	20.0
Endosulfan Sulfate	23.17	23.02-23.34	11.21/05 = 224	Conf.	23.13	11.01/05 = 220	Conf.	1.8
4,4'-DDT	17.35	17.29-17.41	9.73/05 = 195	Conf. B.L.	—	—	—	—
Methoxychlor	34.74	34.52-35.04	4.22/05 = 181	Conf. B.L.	—	—	—	—
Endrin Ketone	31.60	31.33-31.83	17.27/05 = 345 BL	Conf.	31.53	13.71/05 = 274	Conf.	21.0
Tech. Chlordane								
alpha-Chlordane								
gamma-Chlordane								
Toxaphene								
Araclor - 1016								
Araclor - 1221								
Araclor - 1232								
Araclor - 1242								
Araclor - 1248								
Araclor - 1254								
Araclor - 1260								

64 79

* SEE EXHIBIT B, PART 7 a = poor integration caused by carryover ** CONF. = CONFIRMATION (<20% DIFFERENCE)

Project 3-705-050Column SP 2250/2401

PRIORITY POLLUTANT PESTICIDE DATA

RUN DATE 3/30/85ANALYST: BLHP 5870 LOGBOOK PAGE 167,168

(see reverse for linear regression analyses)

Component	RT (Min)	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std Conc (ng/ml)	Area ₃ X 10 ³	Std Conc (ng/ml)	Area ₃ X 10 ³	Std Conc (ng/ml)	Area ₃ X 10 ³	Std Conc (ng/ml)	Area ₃ X 10 ³
1. α-BHC	1.97-2.01	100	314	50	136	25	58.64	10	15.59	5	6.40		
2. γ-BHC	2.56-2.60		273		121		52.50		14.81		6.43		
3. Heptachlor	3.16-3.20		280		123		56.54		18.92		9.22		
4. Aldrin	3.87-3.89		265		108		46.82		14.74		6.96		
5. Hept. Epox.	6.01-6.03		245		106		49.08		18.40		8.18		
6. Dieldrin	9.41-9.49		198		77.8		36.97		12.58		6.13		
7. Endrin	11.51-11.59		142		54.5		26.39		9.62		4.06		
8. p,p'-DDD	14.35-14.41		126		47.5		21.53		7.73		3.59		
9. p,p'-DDT	17.33-17.45		104		37.5		16.14		5.42		2.72		
10. Endrin Ald.	18.77-19.03		202		87.9		41.26		13.81		5.73		
11. Endrin Ketone	31.55-31.79		132		54.2		25.56		9.42		2.44		
12. Methoxychlor	34.76-34.85	↓	11.5	↓	2.14	↓	.90	↓	—	↓	—		
Pesticide Mix		P-30-1		P-30-2		P-30-3		P-30-4		P-30-5			

Project 3-705-050

Column SP 2250/2401

PRIORITY POLLUTANT PESTICIDE DATA

RUN DATE: 3/30/85

ANALYST: BL

HP 5870 LOGBOOK PAGE 167, 168

(see reverse for linear regression analyses)

Component	RT (Min)	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³
1. B-BHC	3.01-3.03	100	^{3.02} 133	50	^{3.02} 66.2	25	^{3.02} 27.88	10	^{3.02} 10.95	5	^{3.02} 4.44		
2. δ-BHC	3.52-3.56		^{3.53} 268		^{3.54} 119		^{3.54} 40.74		^{3.54} 13.59		^{3.53} 4.76		
3. Endosulfan I	7.61-7.67		^{7.64} 236		^{7.64} 111		^{7.64} 45.52		^{7.65} 18.67		^{7.63} 8.47		
4. p,p'-DDE	9.12-9.14		^{9.13} 264		^{9.13} 115		^{9.13} 42.77		^{9.14} 16.72		^{9.13} 6.67		
5. Endosulfan II	14.14-14.24		^{14.19} 215		^{14.19} 101		^{14.19} 40.44		^{14.22} 16.17		^{14.19} 6.66		
6. Endosulfan Sulfate	23.11-23.35		^{23.21} 146		^{23.21} 70.7		^{23.23} 29.63		^{23.30} 13.33		^{23.23} 5.75		
7. Dibutyl Chlorendate	31.65-31.83	(M.S.) ↓	^{31.71} 132	↓	^{31.73} 66.9	↓	^{31.74} 29.25	↓	^{31.79} 12.41	↓	^{31.74} 2.13		
Pesticide Mix			P-31-1		P-31-2		P-31-3		P-31-4		P-31-5		

18 AT

LINEAR REGRESSION RESULTS

SP0250/2401

3/30/85

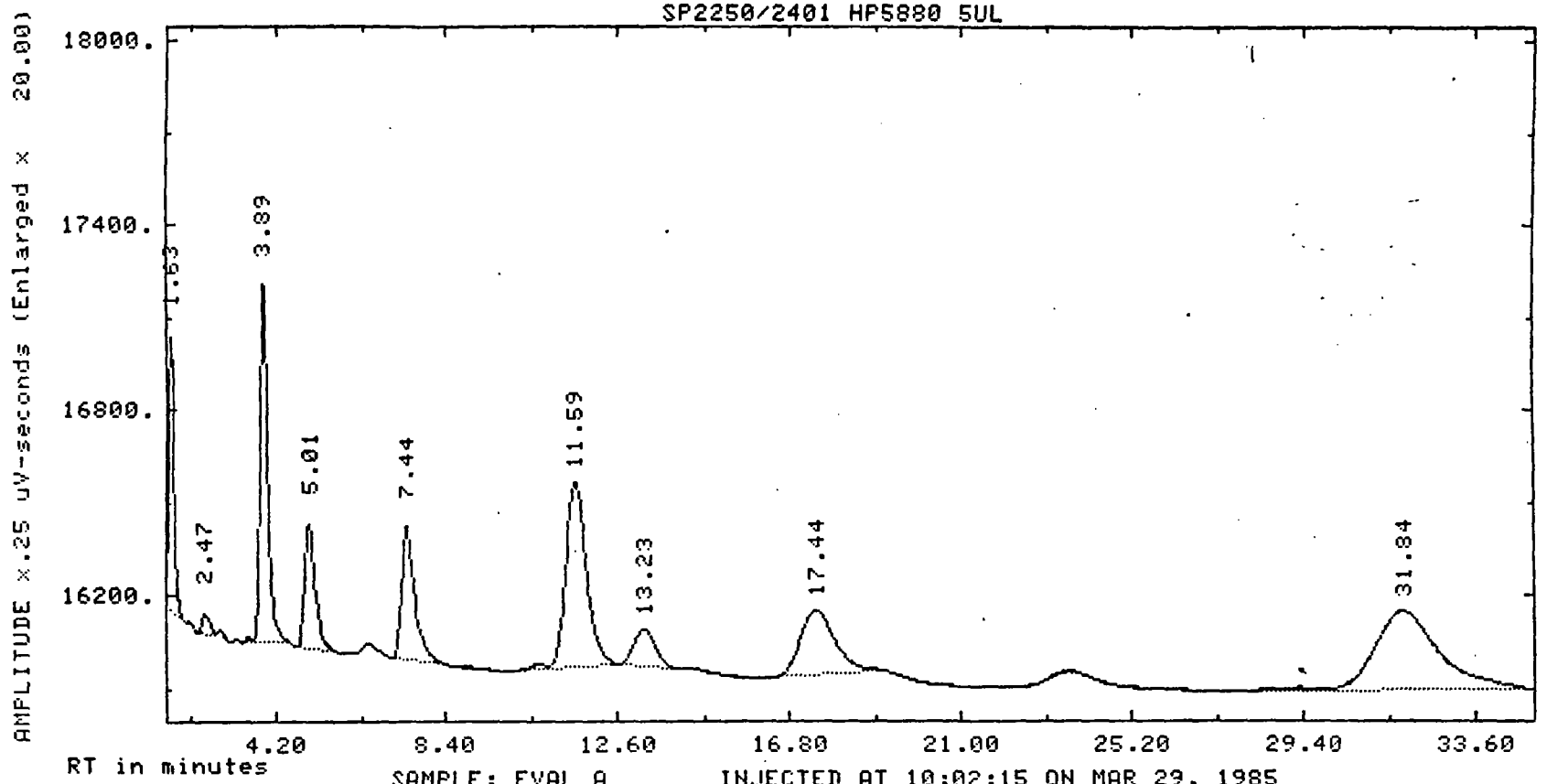
Component	Correlation Coefficient	Y-intercept	Slope
1. β -BHC			
2. δ -BHC			
3. Endosulfan I			
4. p,p'-DDE			
5. Endosulfan II			
6. Endosulfan Sulfate			
7. Dibutyl Chlorendate	.99939	-3.07	1.36

IV
28/82

Case 3969

SP2250/2401 HP5880 5UL

58 II



RT in minutes

SAMPLE: EVAL A INJECTED AT 10:02:15 ON MAR 29, 1985
Method: PRIME Raw: R5001 Proc: P5001

READY
*LI,P,P5001

PROCESSED DATA FILE: P5001 ON CRN 21

Case 3969

APR 2, 1985 11:29:44

REPORT: 97 CHANNEL: 2 # PEAKS: -11 SP2250/2401 HP5080 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.04	0.00	1.00000	77 BV	.146	
2	1.16	0.00	1.00000	159 VB	.300	
3	1.63	0.00	1.00000	3861 BB	7.260	
4	2.47	0.00	1.00000	347 BB	.652	
5	3.89	0.00	1.00000	7451 BB	14.009	Aldrin
6	5.01	0.00	1.00000	3009 BB	6.785	
7	7.44	0.00	1.00000	5256 BB	9.883	
8	11.59	0.00	1.00000	10520 BB	19.779	Endrin
9	13.23	0.00	1.00000	2138 BB	4.020	
10	17.44	0.00	1.00000	5024 BB	10.051	pp'-DDT
11	31.84	0.00	1.00000	13943 BB	26.216	Dibutylchloroendate

TOTAL AREA = 53187 TOTAL AREA % = 100.000

SAMPLE: EVAL A INJECTED AT 10:02:15 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 1 BTL: 1

SL-WDTH MV/MIN DELAY MIN-AR BUNCH REPORT
.500 .300 1.00 50 AUTO MEDIUM

SUP-UNK DVT ID-LVL REF-RTW %RTW %DIL-F ISO
NO 0.00 0 .30 5.0 100.00 NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

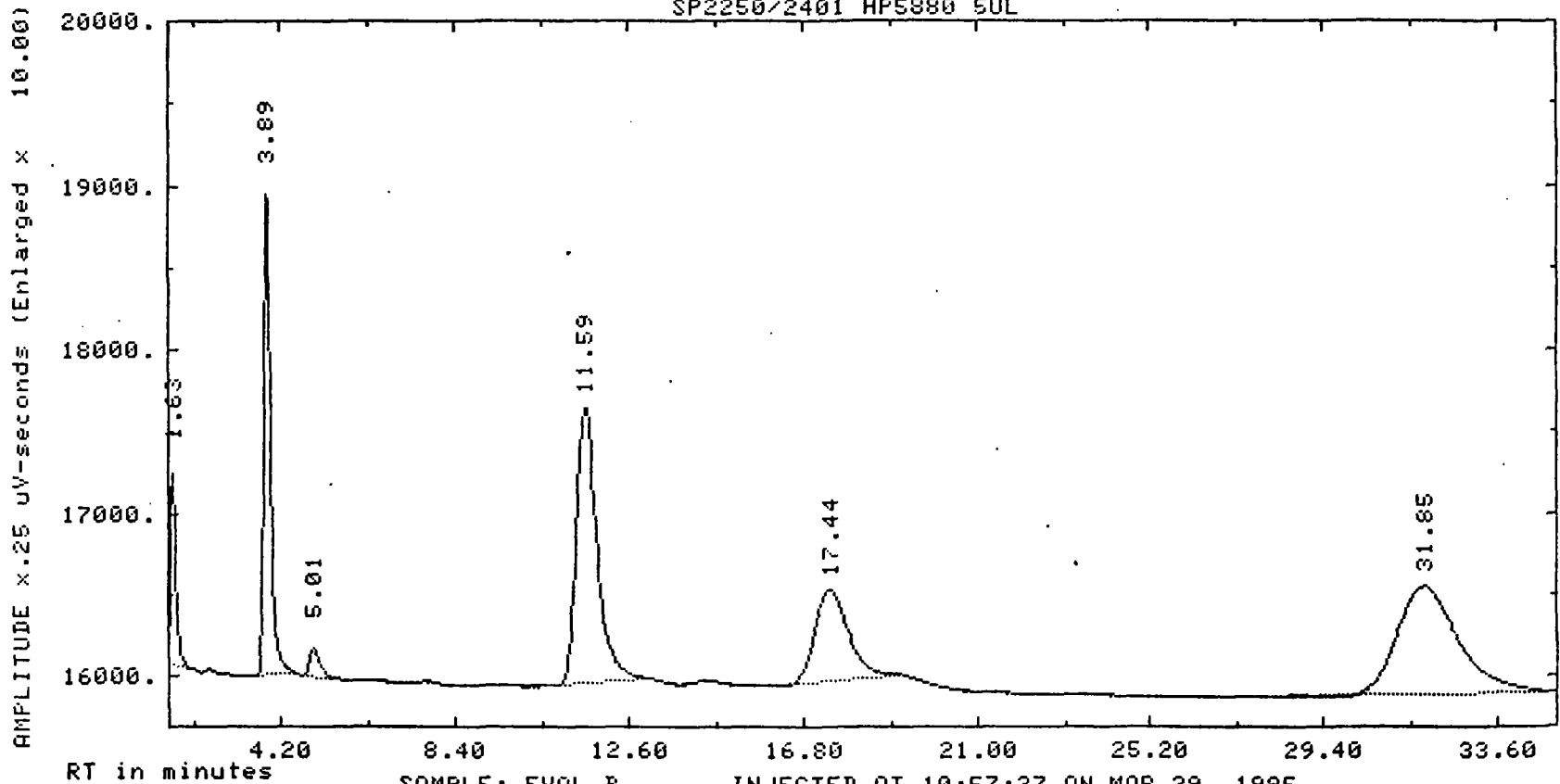
RAW DATA FILE: R5001 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 84

Case 3969

SP2250/2401 HP5880 5UL



SAMPLE: EVAL B INJECTED AT 10:57:27 ON MAR 29, 1985
Method: PRIME Raw: R5002 Proc: P5002

IV 85

APR 2, 1985 11:30:00

REPORT: 93 CHANNEL: 2 # PEAKS: -6 SP2250/2401 HP5860 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	5071 BB	4.882	
2	3.89	0.00	1.00000	17693 BB	17.032	Aldrin
3	5.01	0.00	1.00000	1424 BB	1.371	
4	11.59	0.00	1.00000	29429 BB	28.329	Endrin
5	17.44	0.00	1.00000	15772 BB	15.182	pp'-DDT
6	31.85	0.00	1.00000	34494 BF	33.205	Dibutylchlorodate

TOTAL AREA = 103883 TOTAL AREA % = 100.000

SAMPLE: EVAL B INJECTED AT 10:57:27 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 2 BTL: 2

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISG
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

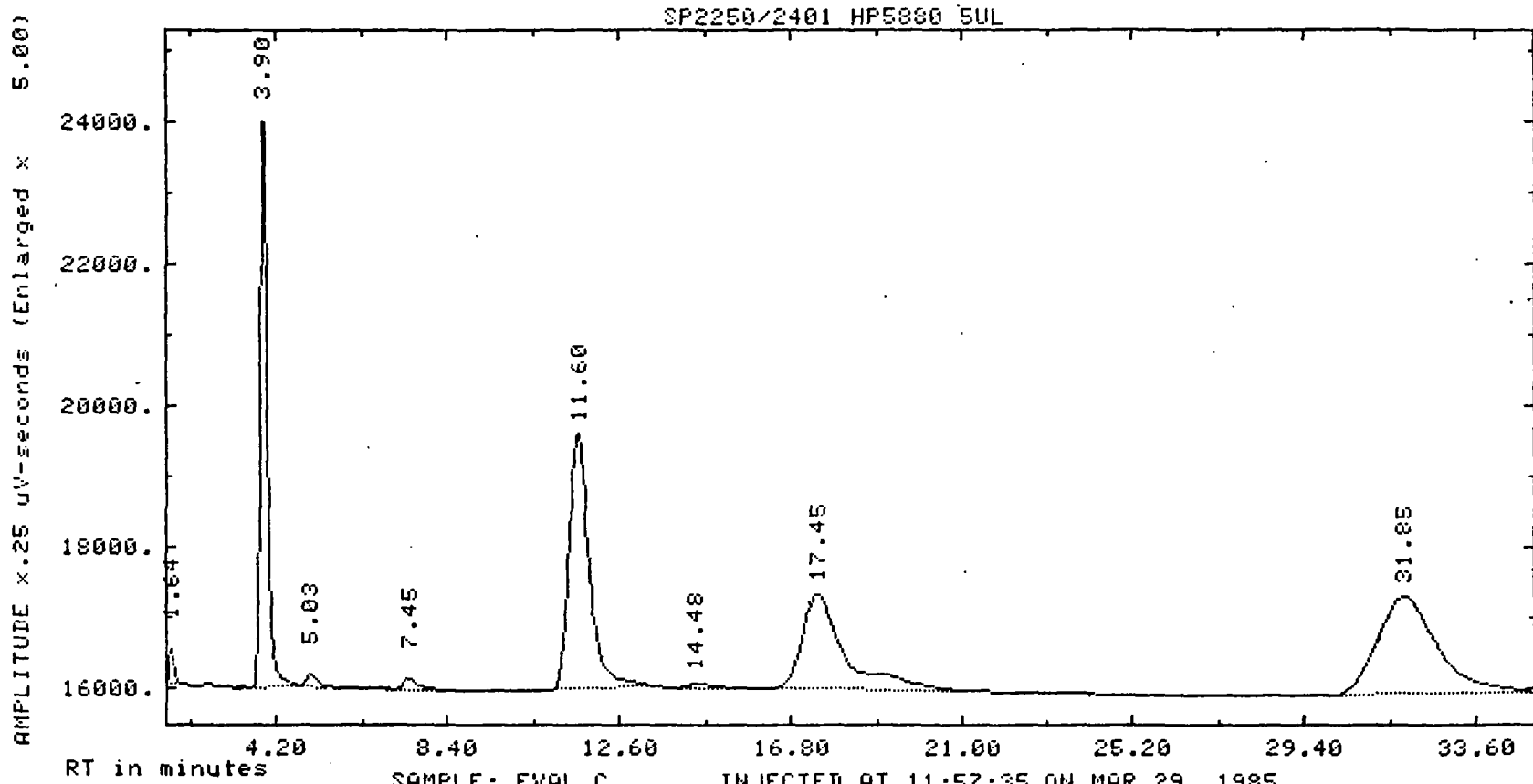
ENDED NOT ON BASELINE

RAW DATA FILE: P5002 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 86

Case 3969



SAMPLE: EVAL C INJECTED AT 11:57:35 ON MAR 29, 1985
Method: PRIME Raw: R5003 Proc: P5003

68 AF

APR 2, 1985 11:30:29

REPORT: 99 CHANNEL: 2 # PEAKS: -8 SP2250/2401 HP5980 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.64	0.00	1.00000	1908 BB	.020	
2	3.90	0.00	1.00000	46303 BB	20.199	Aldrin
3	5.03	0.00	1.00000	1515 BB	.651	
4	7.45	0.00	1.00000	1741 BB	.749	
5	11.60	0.00	1.00000	62034 BB	26.664	Endrin
6	14.40	0.00	1.00000	1004 BB	.432	
7	17.45	0.00	1.00000	47777 BB	20.536	pp'-DDT
8	31.85	0.00	1.00000	69600 BF	29.950	Dibutylchloroendate

TOTAL AREA = 232054 TOTAL AREA % = 100.000

SAMPLE: EVAL C INJECTED AT 11:57:35 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 3 BTL: 3

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	XRTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

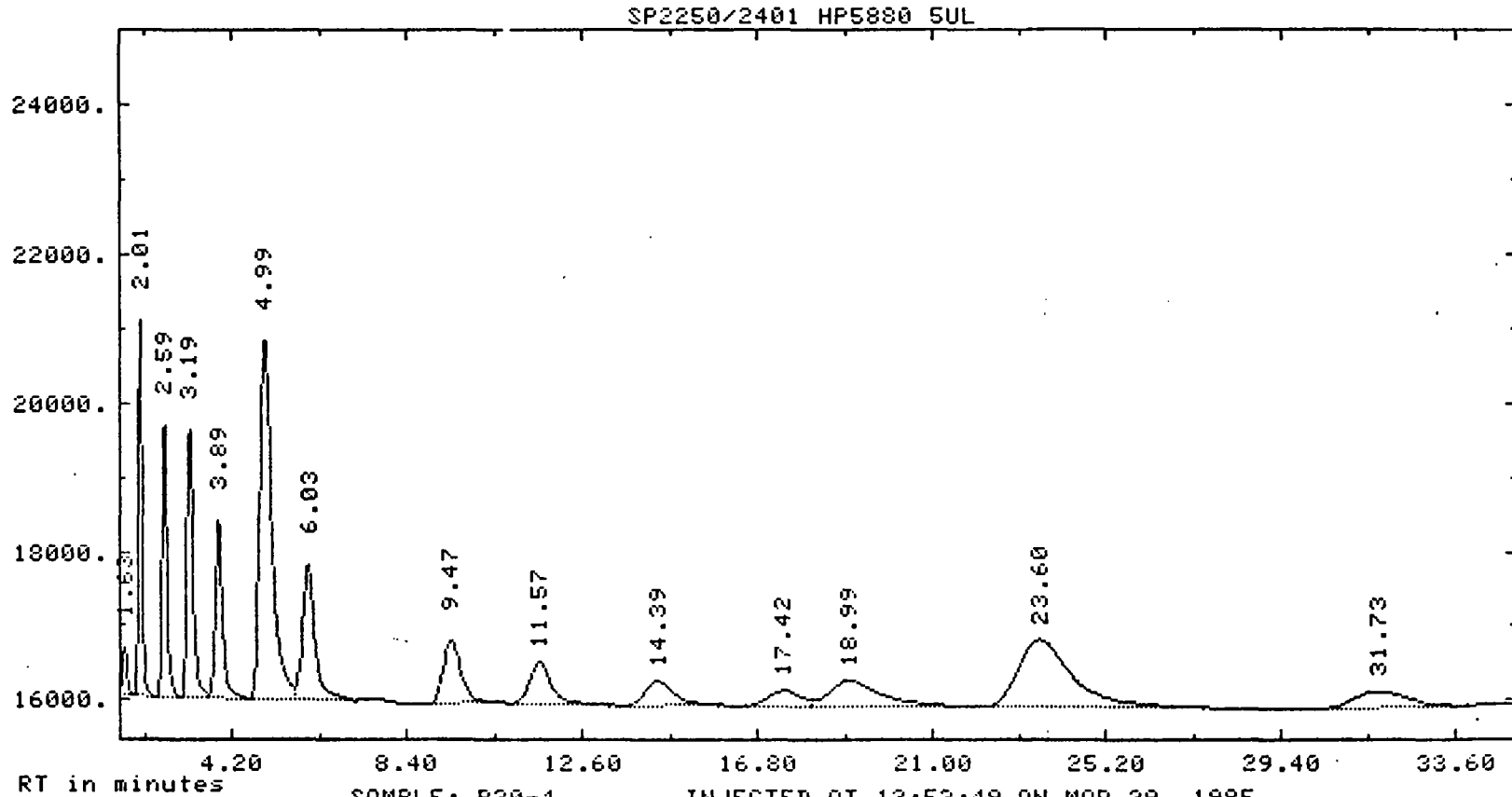
RAW DATA FILE: R5003 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 88

Case 3969

AMPLITUDE x.25 uV-seconds (Enlarged x 150.00)



RT in minutes

SAMPLE: P30-4 INJECTED AT 12:52:49 ON MAR 29, 1985
Method: PRIME Raw: R5004 Proc: P5004

18 II

*LI,P,P5004

PROCESSED DATA FILE: P5004 ON CRN 21

Case 3969

APR 2, 1985 11:30:52

REPORT: 100 CHANNEL: 2 # PEAKS: -14 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	2459 BV	1.048	
2	2.01	0.00	1.00000	15589 VB	6.640	α -BHC
3	2.59	0.00	1.00000	14913 BV	6.310	δ -BHC
4	3.19	0.00	1.00000	10923 VU	8.061	Heptachlor
5	3.89	0.00	1.00000	14785 VB	6.298	Aldrin
6	4.99	0.00	1.00000	50244 BV	21.403	
7	6.03	0.00	1.00000	10395 VB	7.836	Heptachlor Epoxide
8	9.47	0.00	1.00000	12577 BB	5.350	Dieldrin
9	11.57	0.00	1.00000	9618 BB	4.097	Endrin
10	14.39	0.00	1.00000	7734 BB	3.295	pp'-DDO
11	17.42	0.00	1.00000	5421 BV	2.309	pp'-DDT
12	18.99	0.00	1.00000	13808 VB	5.882	Endrin Aldelyde
13	23.60	0.00	1.00000	40969 BB	17.452	
14	31.73	0.00	1.00000	9416 DB	4.011	Endrin Ketene

TOTAL AREA = 234752 TOTAL AREA % = 100.000

SAMPLE: P30-4 INJECTED AT 12:52:49 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 4 BTL: 4

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

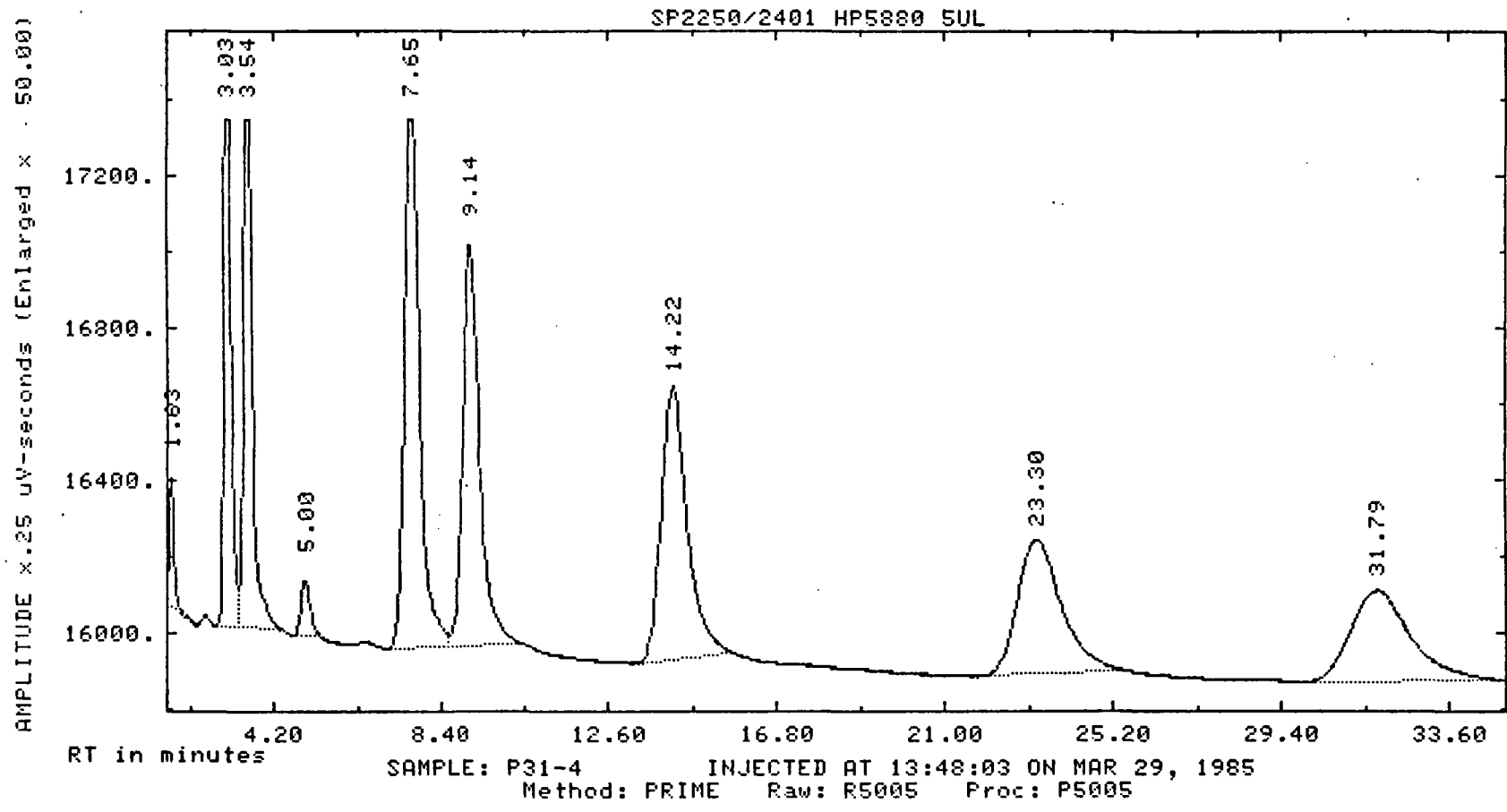
RAW DATA FILE: R5004 PARAM FILES= METHOD: SEQ:

DONE
READY

TV 90

Case 3969

16 μ R



*LI,P,P5005

PROCESSED DATA FILE: P5005 ON CRN 21

Case 3969

APR 2, 1985 11:31:20

REPORT: 101 CHANNEL: 2 # PEAKS: -9 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	1357 BB	1.300	
2	3.03	0.00	1.00000	10952 BV	10.493	B-BHC
3	3.54	0.00	1.00000	13585 VB	13.016	γ-BHC
4	5.00	0.00	1.00000	1182 BB	1.132	
5	7.65	0.00	1.00000	18671 BV	17.888	Endosulfan I
6	9.14	0.00	1.00000	16715 VB	16.014	PP'-DOE
7	14.22	0.00	1.00000	16173 BB	15.495	Endosulfan II
8	23.30	0.00	1.00000	13326 BB	12.767	Endosulfan Sulfate
9	31.79	0.00	1.00000	12414 BB	11.894	D-butylchloroendate

TOTAL AREA = 104374 TOTAL AREA % = 100.000

SAMPLE: P31-4 INJECTED AT 13:48:03 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 5 BTL: 5

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

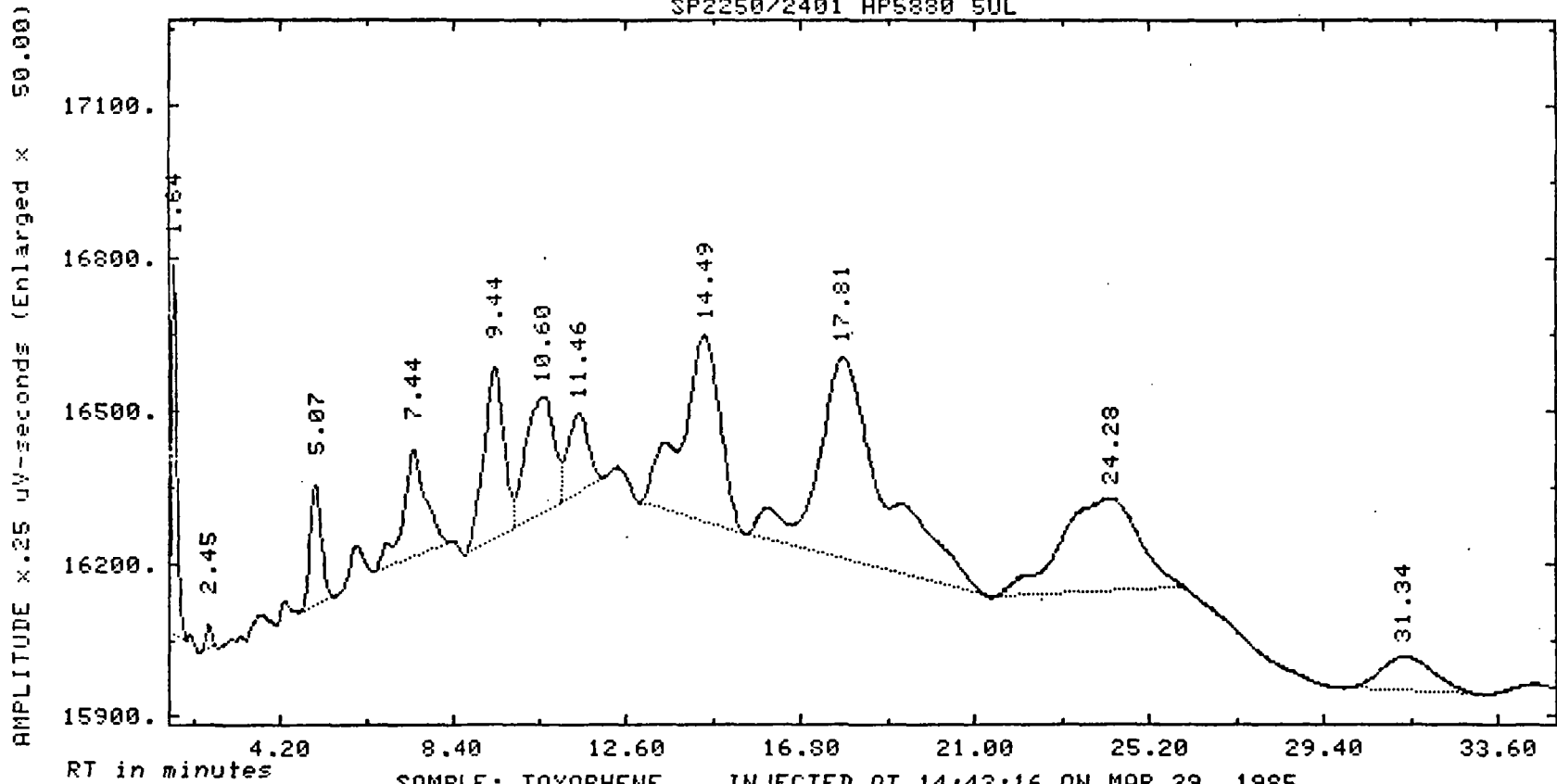
RAW DATA FILE: R5005 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 92

Case 3969

SP2250/2401 HP5880 5UL



SAMPLE: TOXAPHENE INJECTED AT 14:43:16 ON MAR 29, 1985
Method: PRIME Raw: R5006 Proc: P5006

86 III

*LI,P,P5006

PROCESSED DATA FILE: P5006 ON CRN 21

Case 3969

APR 2, 1985 11:31:45

REPORT: 102 CHANNEL: 2 # PEAKS: -12 SP2250/2401 HP5800 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.17	0.00	1.00000	319 BB	.439	
2	1.64	0.00	1.00000	3134 BB	4.309	
3	2.45	0.00	1.00000	251 BB	.345	
4	5.07	0.00	1.00000	2347 BB	3.228	
5	7.44	0.00	1.00000	3931 BB	5.407	
6	9.44	0.00	1.00000	5589 BV	7.607	
7	10.60	0.00	1.00000	5281 VV	7.263	
8	11.46	0.00	1.00000	2661 VD	3.659	
9	14.49	0.00	1.00000	11896 BB	16.360	
10	17.81	0.00	1.00000	22227 BB	30.569	
11	24.28	0.00	1.00000	12084 BB	16.619	
12	31.34	0.00	1.00000	2992 BB	4.114	

TOTAL AREA = 72713 TOTAL AREA % = 100.000

SAMPLE: TOXAPHENE INJECTED AT 14:43:10 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 6 BTL: 6

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

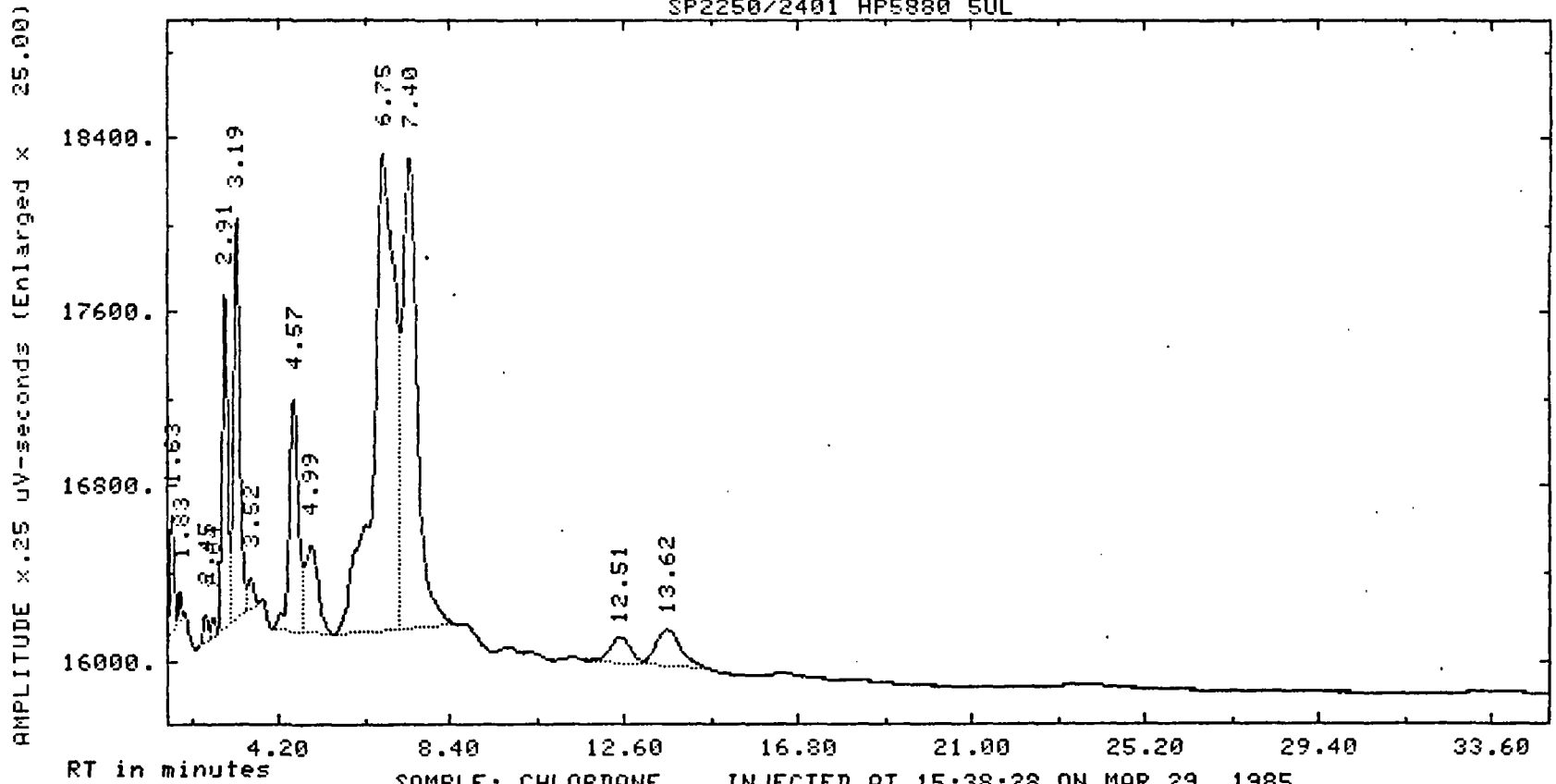
RAW DATA FILE: R5000 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 94

Case 3969

SP2250/2401 HP5880 5UL



56 II

SAMPLE: CHLORDANE INJECTED AT 15:38:28 ON MAR 29, 1985
Method: PRIME Raw: R5007 Proc: P5007

APR 2, 1985 11:32:09

REPORT: 103 CHANNEL: 2 # PEAKS: -15 SP2250/2401 HP5880 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.17	0.00	1.00000	150 BV	.136	
2	1.30	0.00	1.00000	1670 VV	1.518	
3	1.63	0.00	1.00000	1916 VV	1.741	
4	1.83	0.00	1.00000	311 VB	.283	
5	2.45	0.00	1.00000	514 BV	.467	
6	2.64	0.00	1.00000	282 VV	.256	
7	2.91	0.00	1.00000	6934 VV	6.302	
8	3.19	0.00	1.00000	9726 VV	8.840	
9	3.52	0.00	1.00000	683 VB	.621	
10	4.57	0.00	1.00000	6652 BV	7.863	
11	4.99	0.00	1.00000	4029 VB	3.662	
12	6.75	0.00	1.00000	41103 BV	37.356	
13	7.40	0.00	1.00000	28922 VB	26.194	
14	12.51	0.00	1.00000	1981 BV	1.800	
15	13.62	0.00	1.00000	3258 VB	2.951	

TOTAL AREA = 110031 TOTAL AREA % = 100.000

SAMPLE: CHLORDANE INJECTED AT 15:38:28 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 7 BTL: 7

SL-WDTH	NU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	.50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

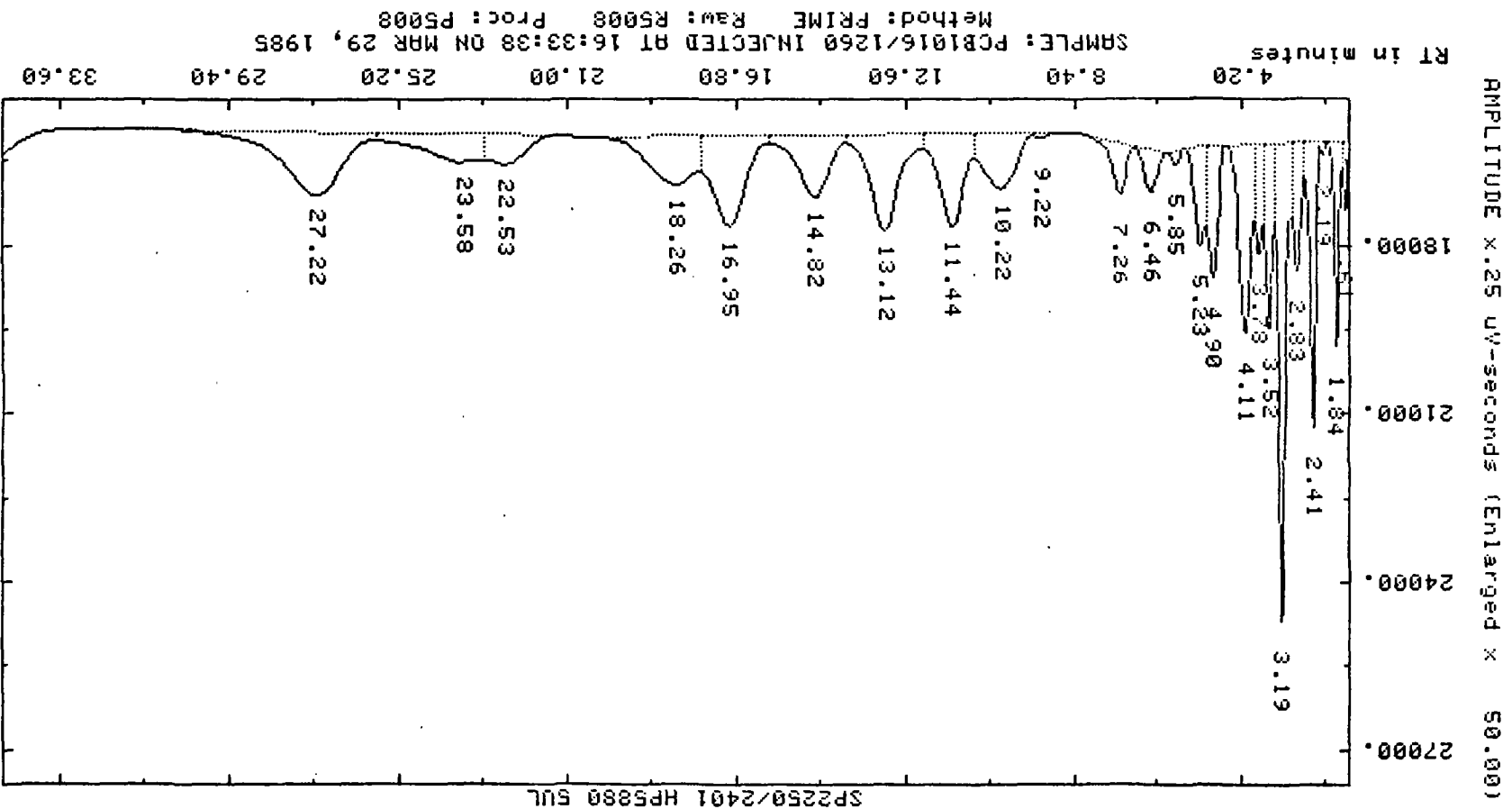
RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: R5007 PARAM FILES= METHOD: SEQ:

DONE
READY

LG AT



Case 3969

APR 2, 1985 11:32:36

REPORT: 104 CHANNEL: 2 # PEAKS: -25 SP2250/2401 HP5980 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.46	0.00	1.00000	1000 BV	.184	
2	1.61	0.00	1.00000	4231 VV	.773	
3	1.84	0.00	1.00000	14369 VV	2.625	
4	2.19	0.00	1.00000	669 VV	.126	
5	2.41	0.00	1.00000	25039 VV	4.575	
6	2.83	0.00	1.00000	14756 VV	2.696	
7	3.19	0.00	1.00000	49950 VV	9.126	
8	3.52	0.00	1.00000	20140 VV	3.690	
9	3.78	0.00	1.00000	10264 VV	1.935	
10	4.11	0.00	1.00000	32667 VV	5.968	
11	4.90	0.00	1.00000	19769 VV	3.612	
12	5.23	0.00	1.00000	15139 VV	2.766	
13	5.85	0.00	1.00000	1473 VB	.269	
14	6.46	0.00	1.00000	8093 BV	1.477	
15	7.26	0.00	1.00000	11789 VB	2.154	
16	9.22	0.00	1.00000	1173 BV	.214	
17	10.22	0.00	1.00000	27756 VV	5.071	
18	11.44	0.00	1.00000	34194 VV	6.247	
19	13.12	0.00	1.00000	43802 VV	8.003	
20	14.02	0.00	1.00000	30802 VV	5.551	
21	16.95	0.00	1.00000	45452 VV	8.304	
22	18.26	0.00	1.00000	33077 VB	6.043	
23	22.53	0.00	1.00000	17414 BV	3.182	
24	23.56	0.00	1.00000	25991 VV	4.749	
25	27.22	0.00	1.00000	57922 VB	10.562	

TOTAL AREA = 547353 TOTAL AREA % = 100.000

SAMPLE: PCB1016/1260 INJECTED AT 16:33:38 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 8 BTL: 8

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

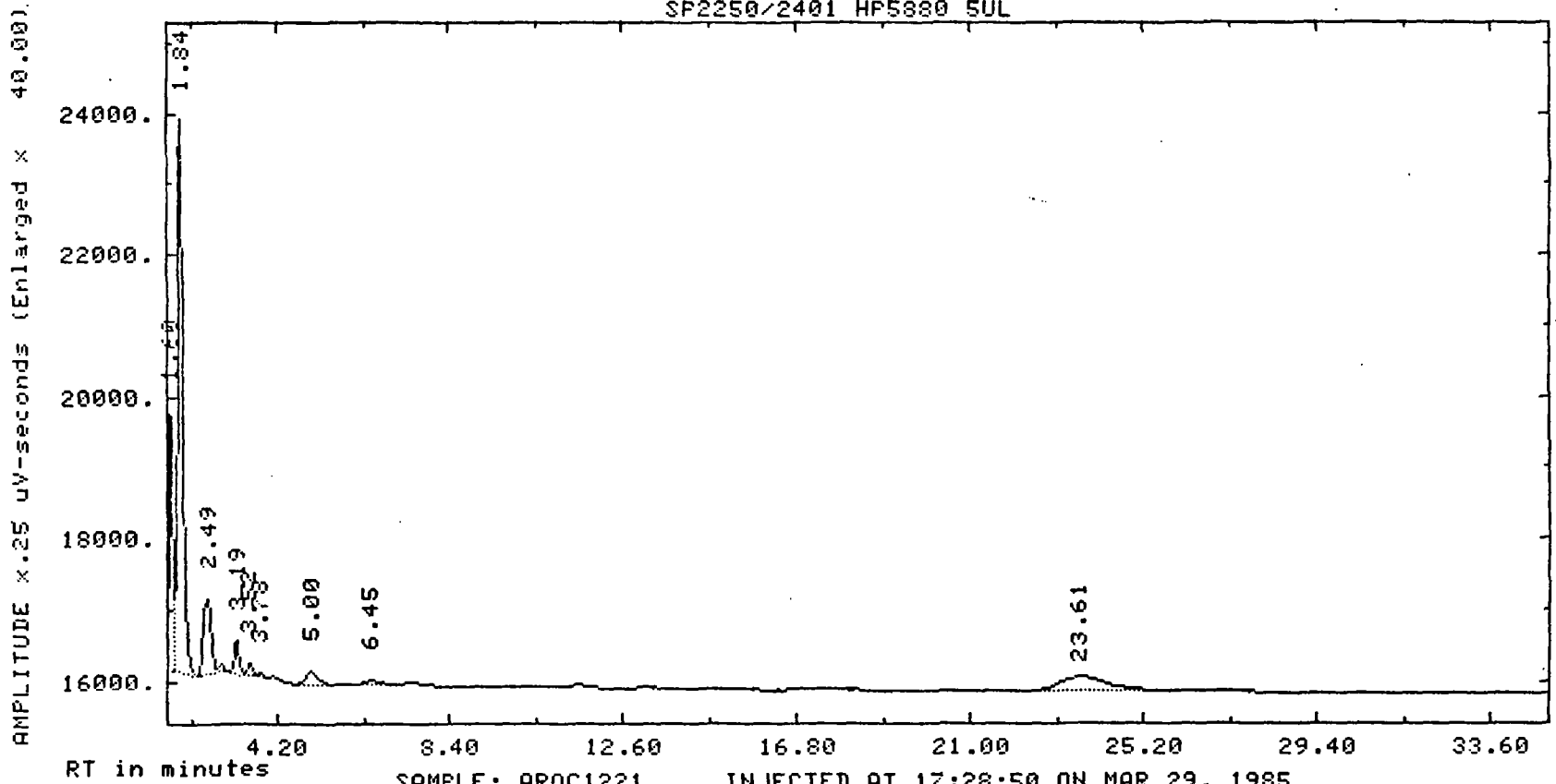
RAW DATA FILE: R5000 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 98

Case 3969

SP2250/2401 HP5880 5UL



AMPLITUDE x.25 uV-seconds (Enlarged x 40.00)

RT in minutes

SAMPLE: AROC1221 INJECTED AT 17:28:50 ON MAR 29, 1985
Method: PRIME Raw: R5009 Proc: P5009

66 II

Case 3969

APR 2, 1985 11:33:10

REPORT: 105 CHANNEL: 2 # PEAKS: -11 SP2250/2401 HP5880 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.27	0.00	1.00000	2340 BV	3.294	
2	1.40	0.00	1.00000	2757 VV	3.881	
3	1.80	0.00	1.00000	11267 VV	15.858	
4	1.84	0.00	1.00000	33238 VV	46.782	
5	2.49	0.00	1.00000	7855 VB	11.057	
6	3.19	0.00	1.00000	2021 BV	2.845	
7	3.52	0.00	1.00000	746 VV	1.050	
8	3.78	0.00	1.00000	225 VB	.317	
9	5.00	0.00	1.00000	1712 BB	2.410	
10	6.45	0.00	1.00000	956 BB	1.346	
11	23.61	0.00	1.00000	7928 BB	11.159	

TOTAL AREA = 71048 TOTAL AREA % = 100.000

SAMPLE: AROC1221 INJECTED AT 17:20:50 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 9 BTL: 9

SL-WIDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

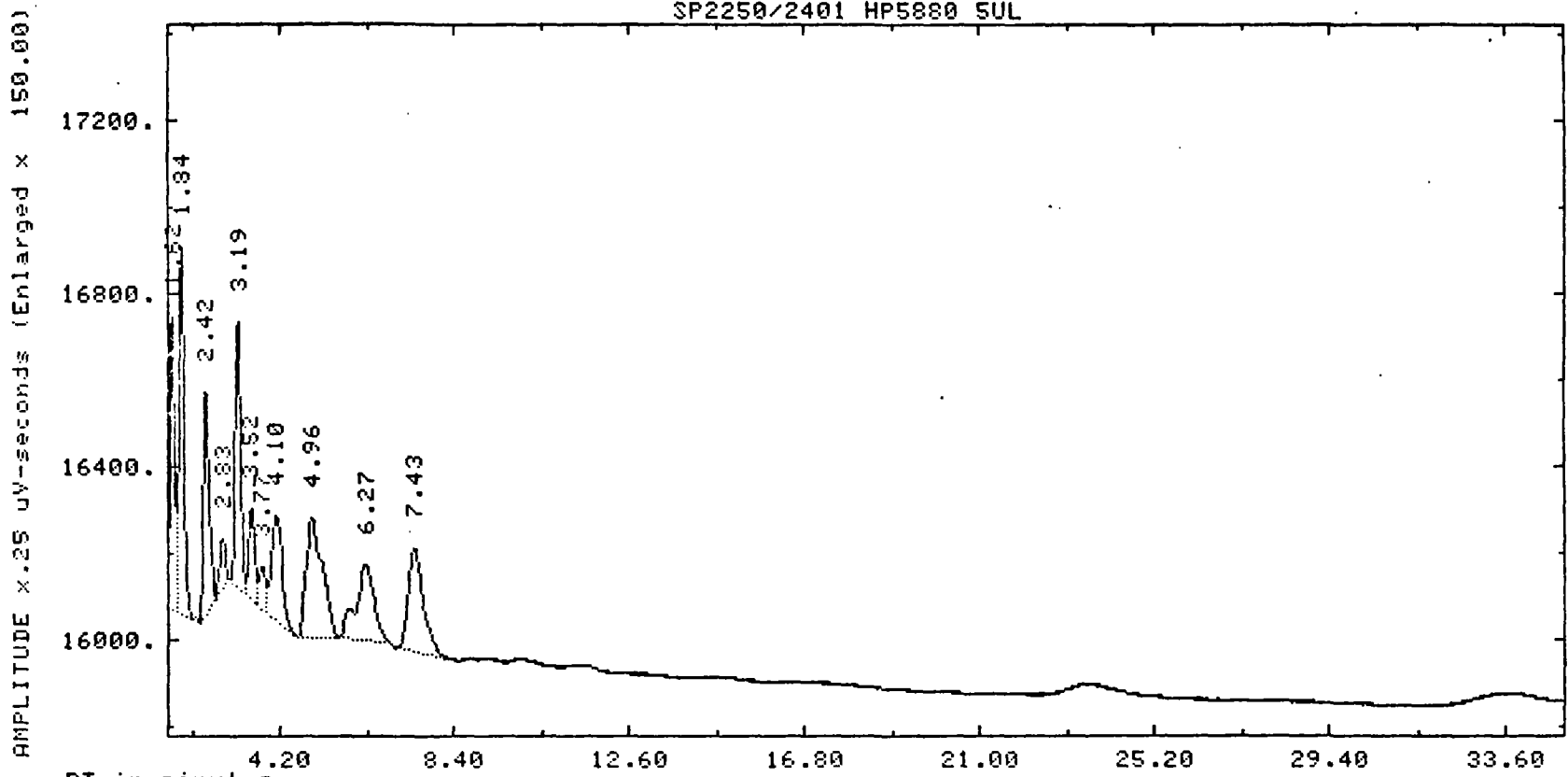
RAW DATA FILE: R5009 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 100

Case 3969

SP2250/2401 HP5880 SUL



AMPLITUDE x.25 uV-seconds (Enlarged x 150.00)

RT in minutes

SAMPLE: AROC1232 INJECTED AT 18:24:03 ON MAR 29, 1985
Method: PRIME Raw: R5010 Proc: P5010

10/11

APR 2, 1985 11:33:34

REPORT: 106 CHANNEL: 2 # PEAKS: -13 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.27	0.00	1.00000	121 BB	.486	
2	1.46	0.00	1.00000	196 BV	.786	
3	1.62	0.00	1.00000	2495 VV	10.008	
4	1.84	0.00	1.00000	3364 VB	13.496	
5	2.42	0.00	1.00000	2478 BV	9.939	
6	2.83	0.00	1.00000	507 VB	2.033	
7	3.19	0.00	1.00000	2868 BV	11.504	
8	3.52	0.00	1.00000	1069 VV	4.289	
9	3.77	0.00	1.00000	489 VV	1.952	
10	4.10	0.00	1.00000	2101 VB	8.430	
11	4.96	0.00	1.00000	3730 BB	14.965	
12	6.27	0.00	1.00000	2593 BB	10.401	
13	7.43	0.00	1.00000	2916 BB	11.699	

TOTAL AREA = 24527 TOTAL AREA % = 100.000

SAMPLE: AROC1232 INJECTED AT 18:24:03 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 10 BTL: 10

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	XRTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.00000 SAMP-AMT: 1.00000

ENDED NOT ON BASELINE

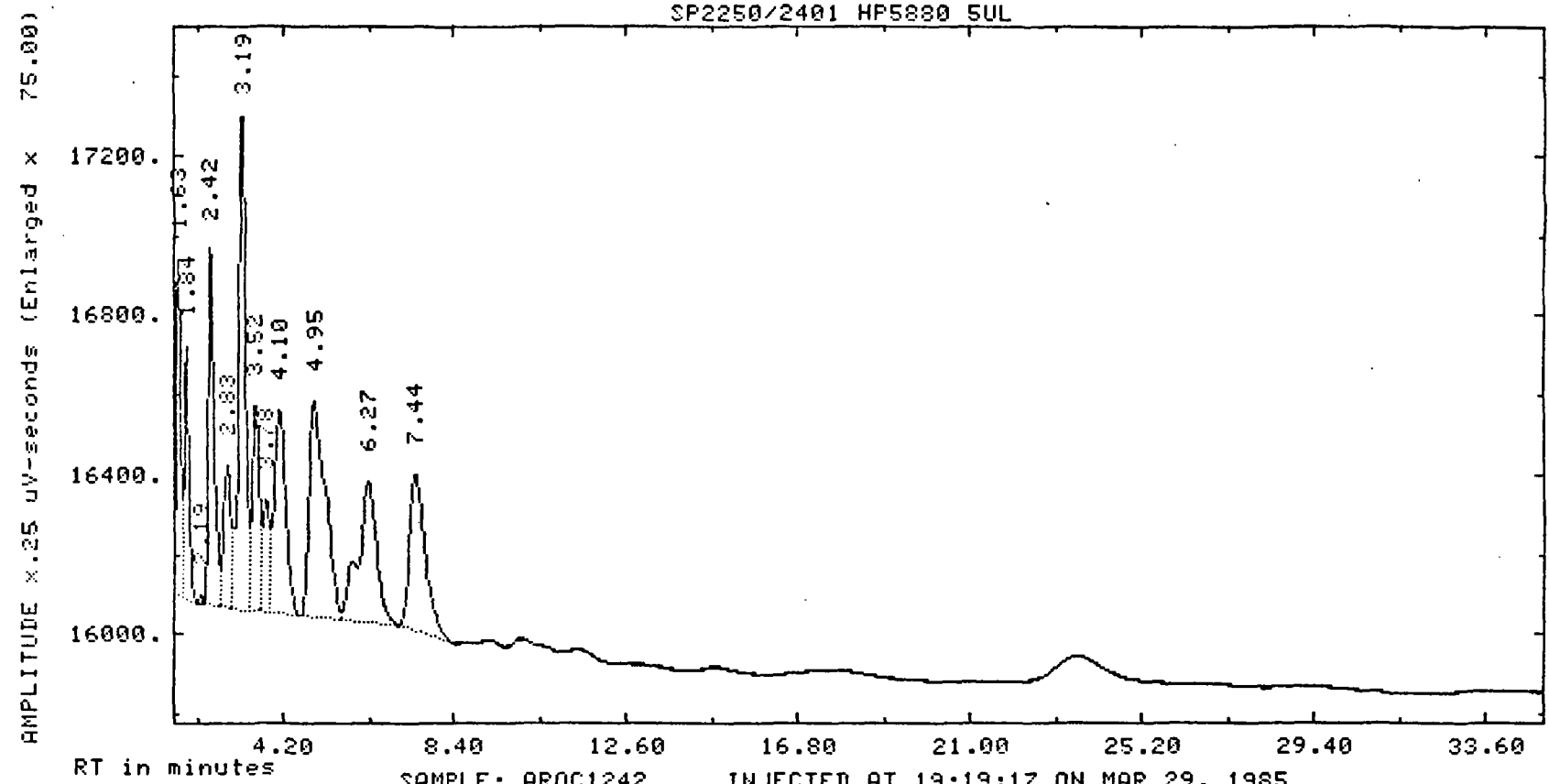
RAW DATA FILE: R5010 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 102

Case 3969

SP2250/2401 HP5880 SUL



SAMPLE: AROC1242 INJECTED AT 19:19:17 ON MAR 29, 1985
Method: PRIME Raw: R5011 Proc: P5011

501 II

*LI,P,P5011

PROCESSED DATA FILE: P5011 ON CRN 21

Case 3969

APR 2, 1985 11:33:59

REPORT: 107 CHANNEL: 2 # PEAKS: -13 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.16	0.00	1.00000	165 BB	.345	
2	1.63	0.00	1.00000	3524 BV	7.344	
3	1.84	0.00	1.00000	2361 VB	4.921	
4	2.19	0.00	1.00000	55 BV	.114	
5	2.42	0.00	1.00000	4322 VV	9.008	
6	2.83	0.00	1.00000	2291 VV	4.775	
7	3.19	0.00	1.00000	7409 VV	15.442	
8	3.52	0.00	1.00000	3137 VV	6.539	
9	3.78	0.00	1.00000	1539 VV	3.207	
10	4.10	0.00	1.00000	4965 VV	10.347	
11	4.95	0.00	1.00000	7440 VV	15.505	
12	6.27	0.00	1.00000	5595 VB	11.661	
13	7.44	0.00	1.00000	5178 BB	10.792	

TOTAL AREA = 47982 TOTAL AREA % = 100.000

SAMPLE: AROC1242 INJECTED AT 19:19:17 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 11 BTL: 11

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%OIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: R5011 PARAM FILES= METHOD: SEQ:

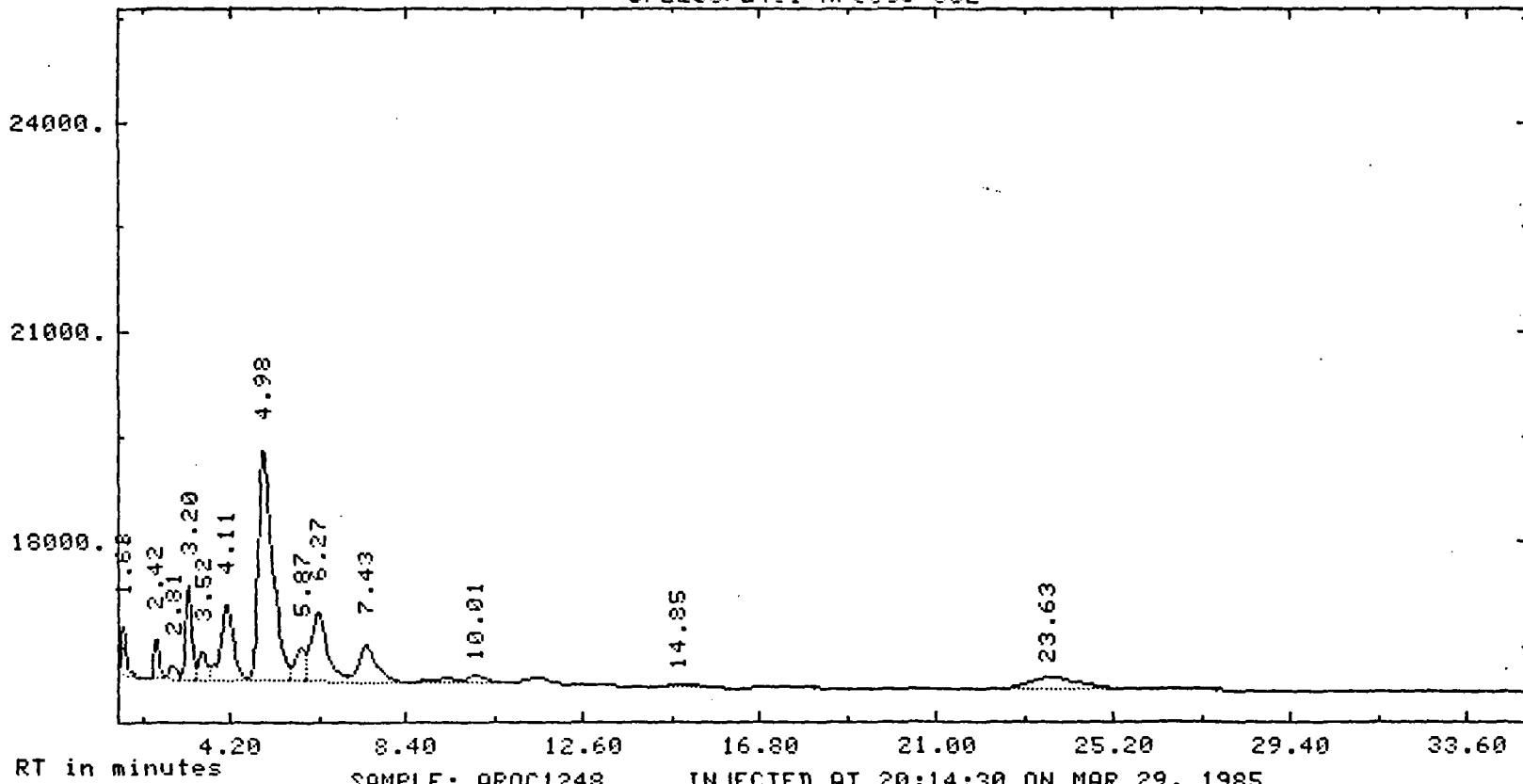
DONE
READY

TV 104

Case 3969

SP2250/2401 HP5880 SUL

AMPLITUDE x.25 uV-seconds (Enlarged x 250.00)



RT in minutes

SAMPLE: AROC1248 INJECTED AT 20:14:30 ON MAR 29, 1985
Method: PRIME Raw: R5012 Proc: P5012

501 II

Case 5167

APR 2, 1985 11:34:32

REPORT: 108 CHANNEL: 2 # PEAKS: -13 SP2250/2401 HPS880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	3128 BB	3.005	
2	2.42	0.00	1.00000	2557 EV	2.456	
3	2.81	0.00	1.00000	1322 VV	1.270	
4	3.20	0.00	1.00000	6995 VV	6.720	
5	3.52	0.00	1.00000	2515 VV	2.417	
6	4.11	0.00	1.00000	11483 VV	11.032	
7	4.98	0.00	1.00000	38834 VV	37.307	
8	5.97	0.00	1.00000	4151 VV	3.988	
9	6.27	0.00	1.00000	13597 VV	13.062	
10	7.43	0.00	1.00000	8394 VB	8.064	
11	10.01	0.00	1.00000	3153 BB	3.030	
12	14.85	0.00	1.00000	1162 BB	1.116	
13	23.63	0.00	1.00000	6801 BB	6.534	

TOTAL AREA = 104093 TOTAL AREA % = 100.000

SAMPLE: AROC1248 INJECTED AT 20:14:30 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 12 BTL: 12

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

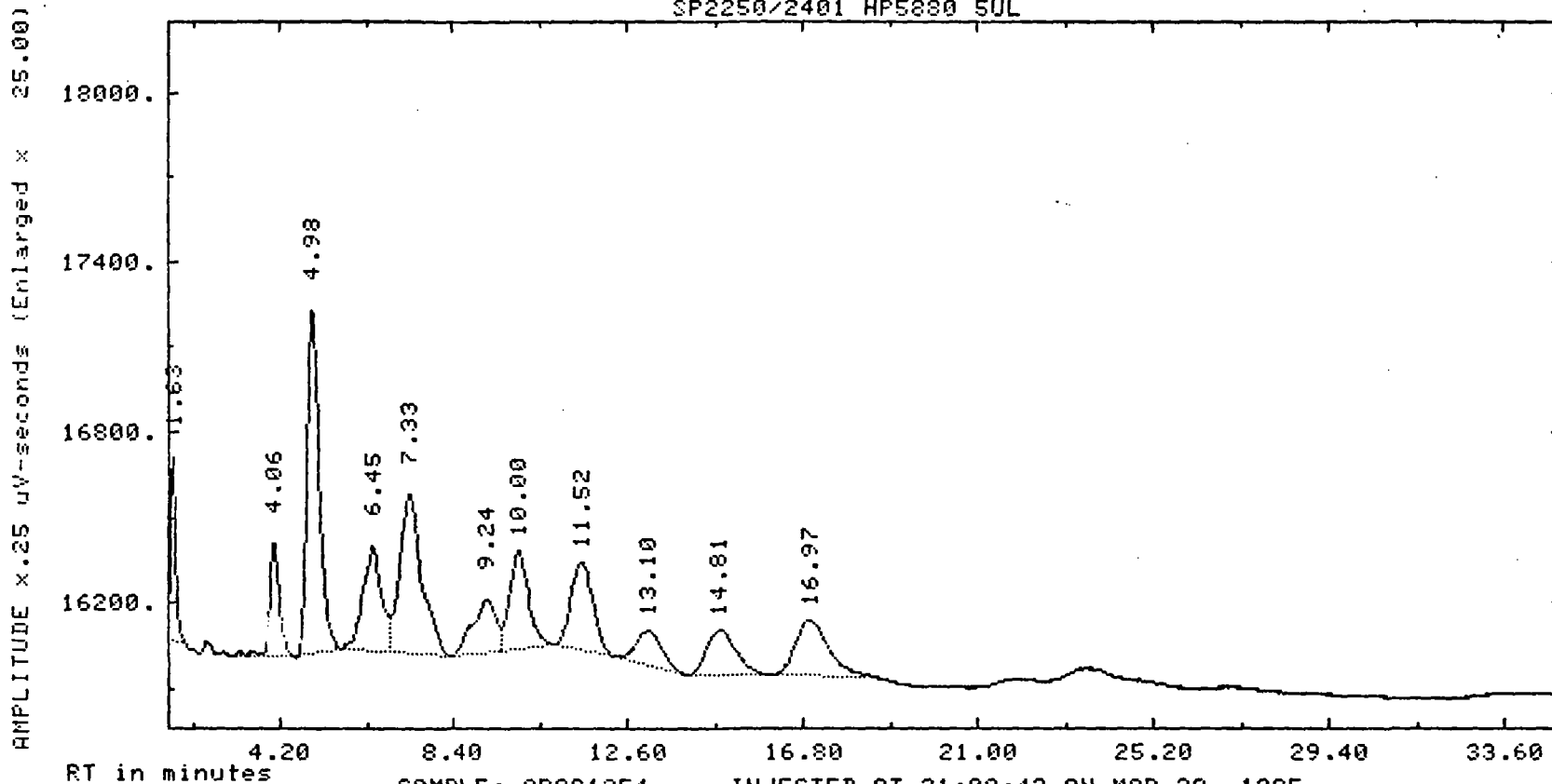
RAW DATA FILE: R5012 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 106

Case 3969

SP2250/2401 HP5880 5UL



AMPLITUDE x.25 μV-seconds (Enlarged x 25.00)

RT in minutes

SAMPLE: AROC1254 INJECTED AT 21:09:42 ON MAR 29, 1985
Method: PRIME Raw: RS013 Proc: P5013

601 II

•LI,P,P5013

PROCESSED DATA FILE: P5013 ON CRN 21

Case 3969

APR 2, 1965 11:34:58

REPORT: 109 CHANNEL: 2 # PEAKS: -11 SP2250/2401 HP5680 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	2647 BB	4.402	
2	4.05	0.00	1.00000	3273 BB	5.442	
3	4.98	0.00	1.00000	12043 BB	20.026	
4	6.45	0.00	1.00000	5792 BV	9.631	
5	7.33	0.00	1.00000	10429 VV	17.341	
6	9.24	0.00	1.00000	3515 VV	5.844	
7	10.00	0.00	1.00000	5434 VB	9.036	
8	11.52	0.00	1.00000	5580 BB	9.279	
9	13.10	0.00	1.00000	2484 BB	4.131	
10	14.81	0.00	1.00000	3727 BB	6.197	
11	16.97	0.00	1.00000	5215 BB	8.671	

TOTAL AREA = 60140 TOTAL AREA % = 100.000

SAMPLE: AROC1254 INJECTED AT 21:08:42 ON MAR 23, 1965

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 13 BTL: 13

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: R5013 PARAM-FILES= METHOD: SEQ:

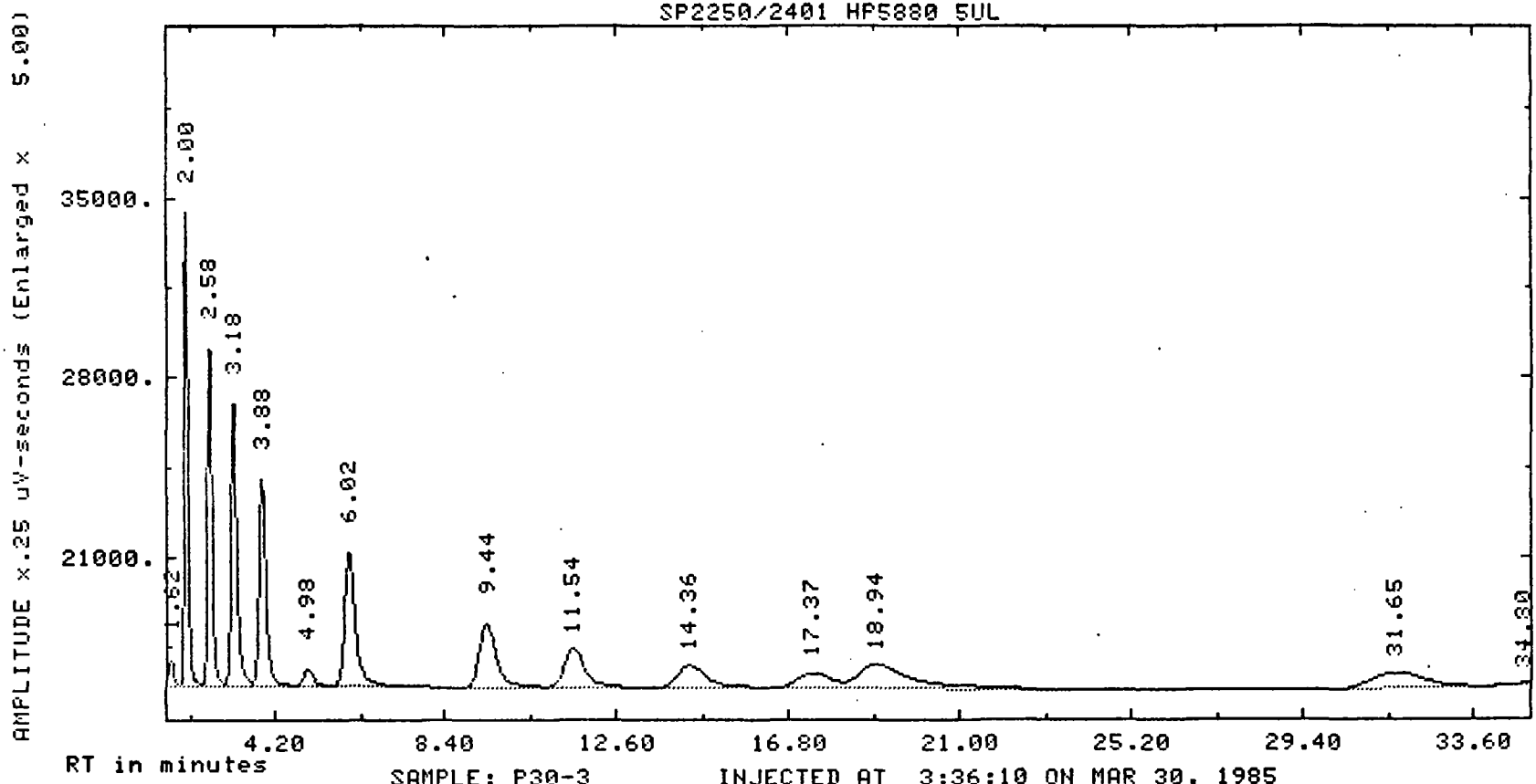
DONE
READY

IV 108

Case 3969

SP2250/2401 HP5880 5UL

601 II



SAMPLE: P30-3 INJECTED AT 3:36:10 ON MAR 30, 1985
Method: PRIME Raw: RS020 Proc: P5020

APR 2, 1985 11:36:17

REPORT: 116 CHANNEL: 2 # PEAKS: -15 SP2250/2401 HP5680 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.17	0.00	1.00000	124 BB	.028	
2	1.62	0.00	1.00000	4159 BV	.940	
3	2.00	0.00	1.00000	58641 UV	13.249	α -BHC
4	2.58	0.00	1.00000	52501 UV	11.862	δ -BHC
5	3.18	0.00	1.00000	56591 UV	12.786	Heptachlor
6	3.88	0.00	1.00000	46821 VB	10.579	Aldrin
7	4.98	0.00	1.00000	5937 BB	1.341	
8	6.02	0.00	1.00000	49083 BB	11.090	Heptachlor Epoxide
9	9.44	0.00	1.00000	36966 BV	8.352	Dieldrin
10	11.54	0.00	1.00000	26393 VB	5.963	Endrin
11	14.36	0.00	1.00000	21529 BB	4.864	pp'-DDD
12	17.37	0.00	1.00000	16141 BV	3.647	pp'-DDT
13	18.94	0.00	1.00000	41256 VB	9.321	Endrin Aldehyde
14	31.65	0.00	1.00000	25555 BV	5.774	Endrin Ketone
15	34.80	0.00	1.00000	897 VF	.203	Methoxychlor

TOTAL AREA = 442596 TOTAL AREA % = 100.000

SAMPLE: P30-3 INJECTED AT 3:36:10 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 20 BTL: 20

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.008 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.00000 SAMP-AMT: 1.00000

ENDED NOT ON BASELINE

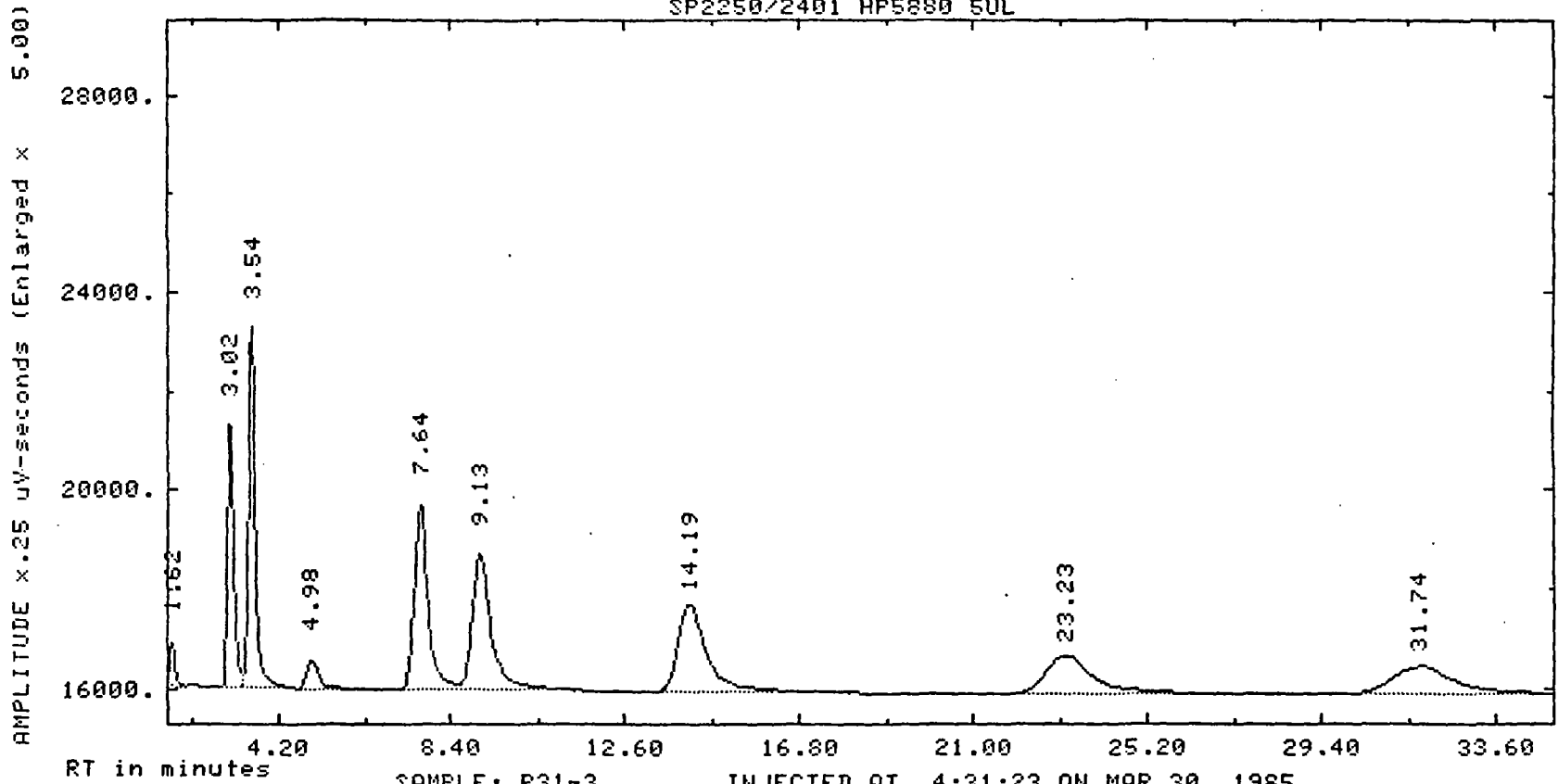
RAW DATA FILE: R5020 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 110

Case 3969

SP2250/2401 HP5880 5UL



SAMPLE: P31-3 INJECTED AT 4:31:23 ON MAR 30, 1985
Method: PRIME Raw: R5021 Proc: P5021

111 II

APR 2, 1985 11:36:45

REPORT: 117 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.19	0.00	1.00000	161 BB	.061	
2	1.62	0.00	1.00000	3486 BB	1.315	
3	3.02	0.00	1.00000	27876 BV	10.513	β-BHC
4	3.54	0.00	1.00000	40736 VB	15.362	δ-BHC
5	4.98	0.00	1.00000	5087 BB	1.919	
6	7.64	0.00	1.00000	45521 BV	17.157	Endosulfan I
7	9.13	0.00	1.00000	42974 VB	16.205	pp'-DDE
8	14.19	0.00	1.00000	40440 BB	15.251	Endosulfan II
9	23.23	0.00	1.00000	29630 BB	11.174	Endosulfan Sulfate
10	31.74	0.00	1.00000	29253 BF	11.032	Dibutylchloride etc

TOTAL AREA = 265164 TOTAL AREA % = 100.000

SAMPLE: P31-3 INJECTED AT 4:31:23 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: 5M03 SUBSQ/SAMP: 1 / 21 BTL: 21

SL-WIDTH MV/MIN DELAY MIN-AR BUNCH REPORT
 .500 .300 1.00 50 AUTO MEDIUM

SUP-UNK DVT ID-LVL REF-RTW %RTW %DIL-F ISO
 NO 0.00 0 .30 5.0 100.00 NO

ACTUAL RUN TIME: 40.008 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

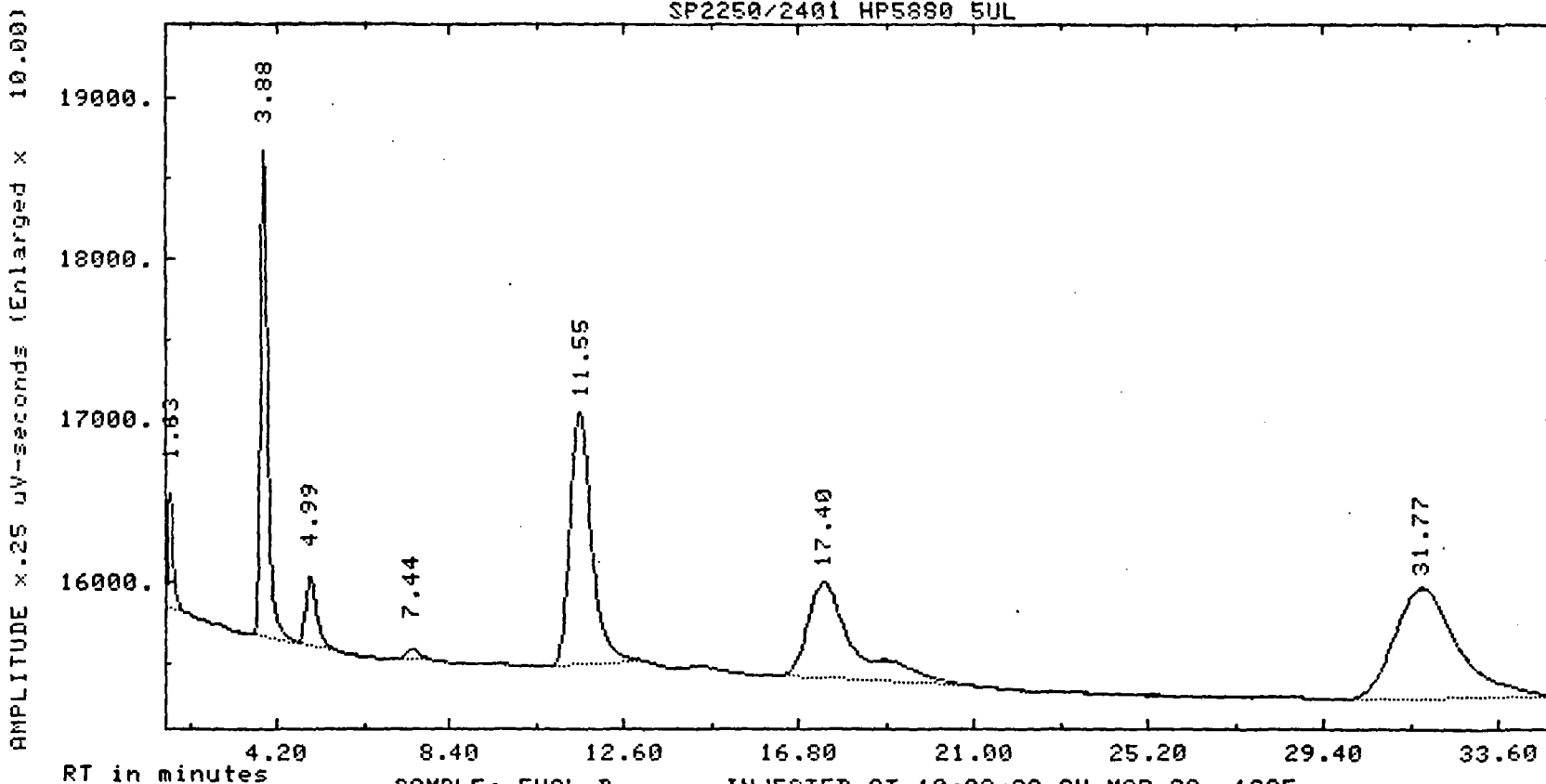
RAW DATA FILE: R5021 PARAM FILES= METHOD: SEQ:

DONE
 READY

IV 112

Case 3969

SP2250/2401 HP5880 5UL



SAMPLE: EVAL B INJECTED AT 10:02:38 ON MAR 30, 1985
Method: PRIME Raw: R5027 Proc: P5027

811 II

*LI,P,P5027

Case 3969

PROCESSED DATA FILE: P5027 ON CRN 21

APR 2, 1985 11:41:41

REPORT: 123 CHANNEL: 2 # PEAKS: -7 SP2250/2401 HPS880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	2897 BB	2.588	
2	3.88	0.00	1.00000	19264 BB	18.319	Aldrin
3	4.99	0.00	1.00000	3858 BB	3.447	
4	7.44	0.00	1.00000	623 BB	.539	
5	11.55	0.00	1.00000	27537 BB	24.693	Endrin
6	17.40	0.00	1.00000	22413 BB	20.026	pp'-DPT
7	31.77	0.00	1.00000	35248 BF	32.387	Dibutylchloroate

TOTAL AREA = 111920 TOTAL AREA % = 100.000

SAMPLE: EVAL B INJECTED AT 10:02:38 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 27 BTL: 27

SL-WDTH	MOV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.008 MINUTES

RF-UNK: 1.00000 STD-AMT: . 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

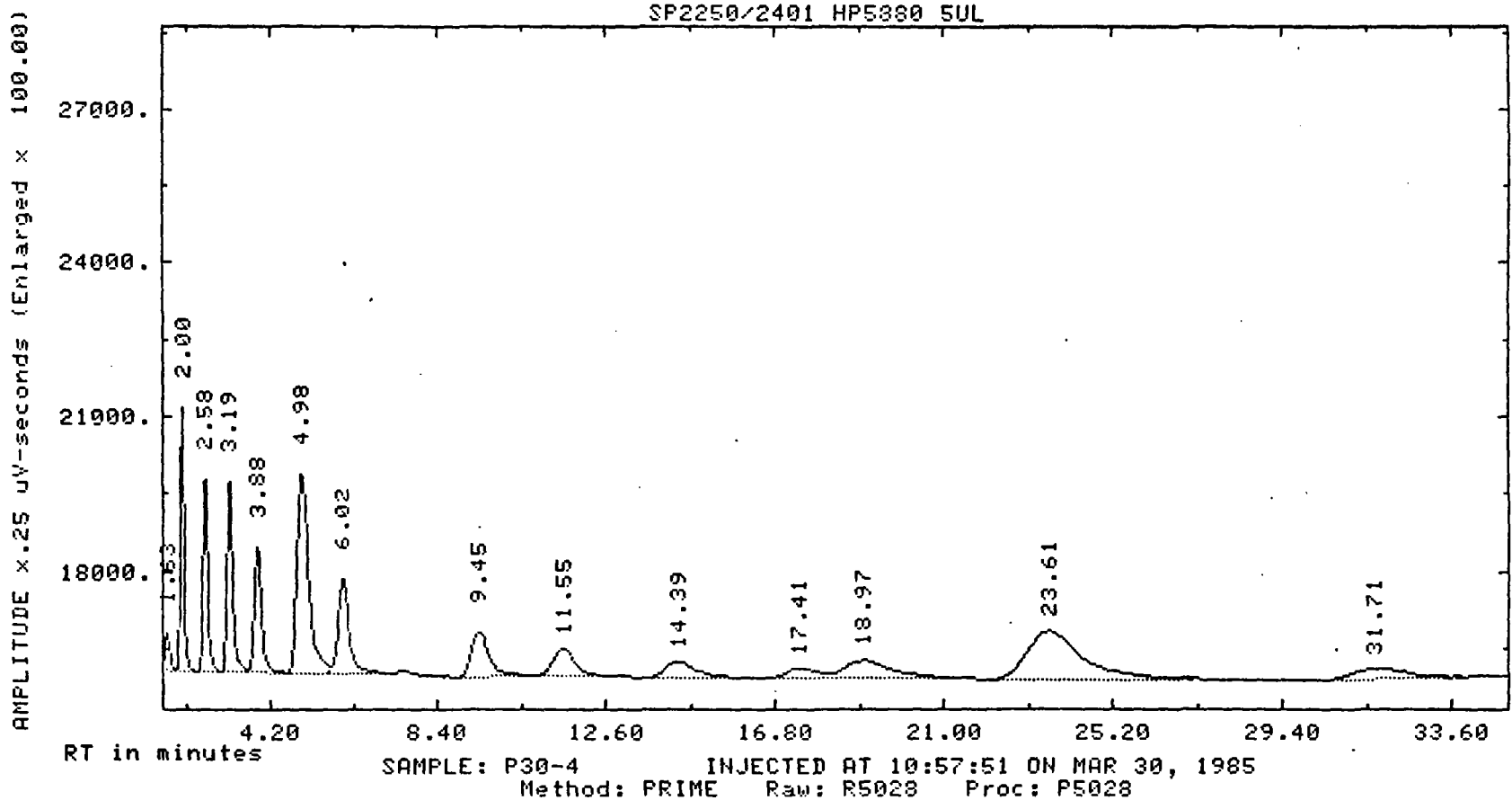
RAW DATA FILE: R5027 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 114

Case 3969

S11 II



*LI,P,P5028

PROCESSED DATA FILE: P5028 ON CRN 21

Case 3969

APR 2, 1985 11:42:03

REPORT: 124 CHANNEL: 2 # PEAKS: -14 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	2942 BV	1.293	
2	2.00	0.00	1.00000	15625 VB	6.866	α -BHC
3	2.58	0.90	1.00000	14940 BV	6.565	β -BHC
4	3.19	0.00	1.00000	19275 VV	8.470	Heptachlor
5	3.88	0.00	1.00000	15071 VB	6.622	Aldrin
6	4.98	0.00	1.00000	39524 BV	17.368	
7	6.02	0.00	1.00000	18397 VB	8.084	Heptachlor Epoxide
8	9.45	0.00	1.00000	12649 BB	5.558	Dieldrin
9	11.55	0.00	1.00000	9572 BB	4.206	Endrin
10	14.39	0.00	1.00000	7585 BB	3.333	PP'-DDD
11	17.41	0.00	1.00000	5538 BV	2.434	PP'-DDT
12	18.97	0.00	1.00000	14146 VB	6.216	Endrin Aldehyde
13	23.61	0.00	1.00000	43489 BB	19.110	
14	31.71	0.00	1.00000	6919 BB	3.075	Endrin Ketone

TOTAL AREA = 227575 TOTAL AREA % = 100.000

SAMPLE: P30-4 INJECTED AT 10:57:51 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 28 STL: 28

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

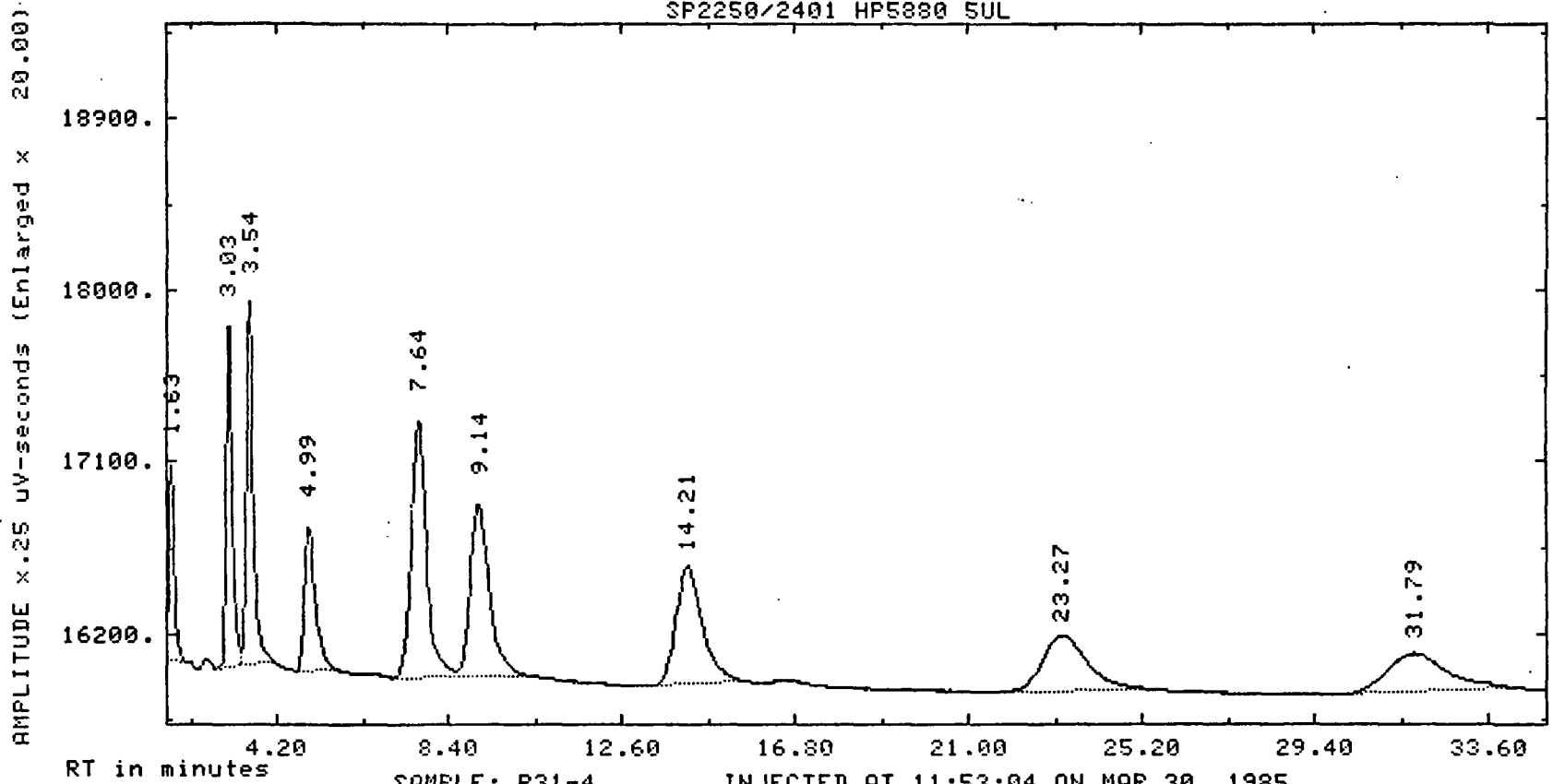
RAW DATA FILE: R5028 PARAM FILES- METHOD: SEQ:

DONE
READY

IV 116

Case 3969

SP2250/2401 HP5880 5UL



AMPLITUDE x.25 uV-seconds (Enlarged x 20.00)

RT in minutes

SAMPLE: P31-4

INJECTED AT 11:53:04 ON MAR 30, 1985

Method: PRIME Raw: R5029 Proc: P5029

L11 AF

*LI,P,P5029

Case 3969

PROCESSED DATA FILE: P5029 ON CRN 21

APR 2, 1985 11:42:28

REPORT: 125 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HP5890 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.18	0.00	1.00000	244 BB	.246	
2	1.63	0.00	1.00000	4559 BB	4.598	
3	3.03	0.00	1.00000	9380 BV	9.459	B-BHC
4	3.54	0.00	1.00000	11050 VB	11.142	g-BHC
5	4.99	0.00	1.00000	7125 BB	7.195	
6	7.64	0.00	1.00000	17528 BV	17.776	Endosulfan I
7	9.14	0.00	1.00000	14432 VB	14.553	pp'-DDE
8	14.21	0.00	1.00000	13892 BB	14.008	Endosulfan II
9	23.27	0.00	1.00000	10735 BB	10.825	Endosulfan Sulfate
10	31.79	0.00	1.00000	10120 BB	10.206	Di-b.tylchlorexate

TOTAL AREA = 99167 TOTAL AREA % = 100.000

SAMPLE: P31-4 INJECTED AT 11:53:04 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 29 BTL: 29

SL-WDTH	MV/NIN	DELAY	NIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.025 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

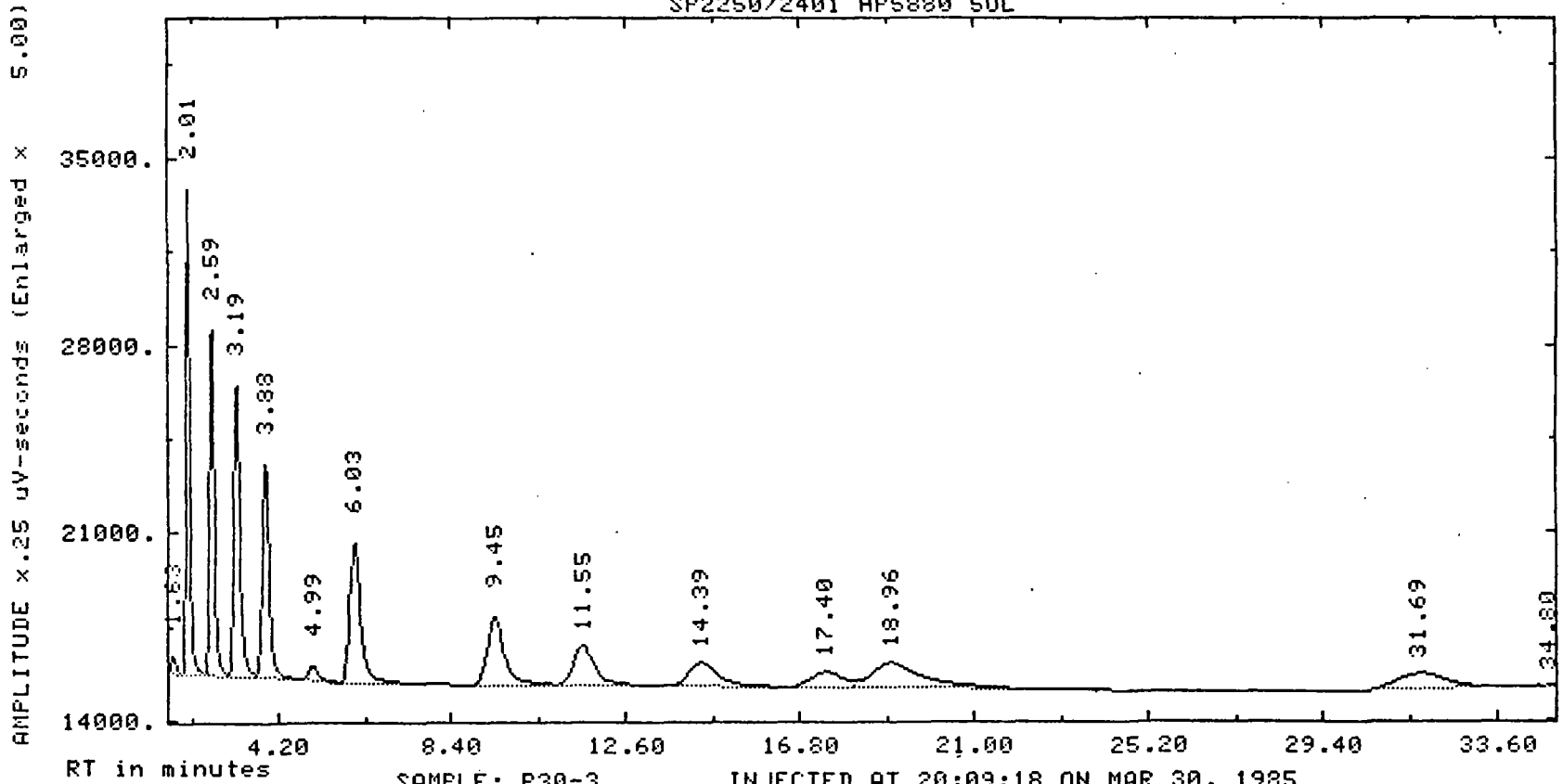
RAW DATA FILE: R5029 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 118

Case 3969

SP2250/2401 HP5880 5UL



SAMPLE: P30-3 INJECTED AT 20:09:18 ON MAR 30, 1985
Method: PRIME Raw: R5038 Proc: P5038

b11 AT

*LI,P,P5038

Case 3969

PROCESSED DATA FILE: P5038 ON CRN 21

APR 2, 1985 11:44:42

REPORT: 134 CHANNEL: 2 # PEAKS: -14 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	2694 BV	.609	
2	2.01	0.00	1.00000	59135 VV	13.370	α -BHC
3	2.59	0.00	1.00000	52781 VV	11.933	γ -BHC
4	3.19	0.00	1.00000	56352 VV	12.740	Heptachlor
5	3.88	0.00	1.00000	46994 VB	10.625	Aldrin
6	4.99	0.00	1.00000	4682 EB	1.059	
7	6.03	0.00	1.00000	48765 EB	11.025	Heptachlor Epoxide
8	9.45	0.00	1.00000	37530 BV	8.485	Dieldrin
9	11.55	0.00	1.00000	26867 VB	6.074	Endrin
10	14.39	0.00	1.00000	22056 EB	4.997	pp'-DDD
11	17.40	0.00	1.00000	16438 BV	3.718	pp'-DDT
12	18.96	0.00	1.00000	41799 VB	9.450	Endrin Aldehyde
13	31.69	0.00	1.00000	25533 BV	5.773	Endrin Ketone
14	34.80	0.00	1.00000	685 VF	.155	Methoxy chlor

TOTAL AREA = 442312 TOTAL AREA % = 100.000

SAMPLE: P30-3 INJECTED AT 20:09:18 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 38 BTL: 28

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	XRTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.008 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

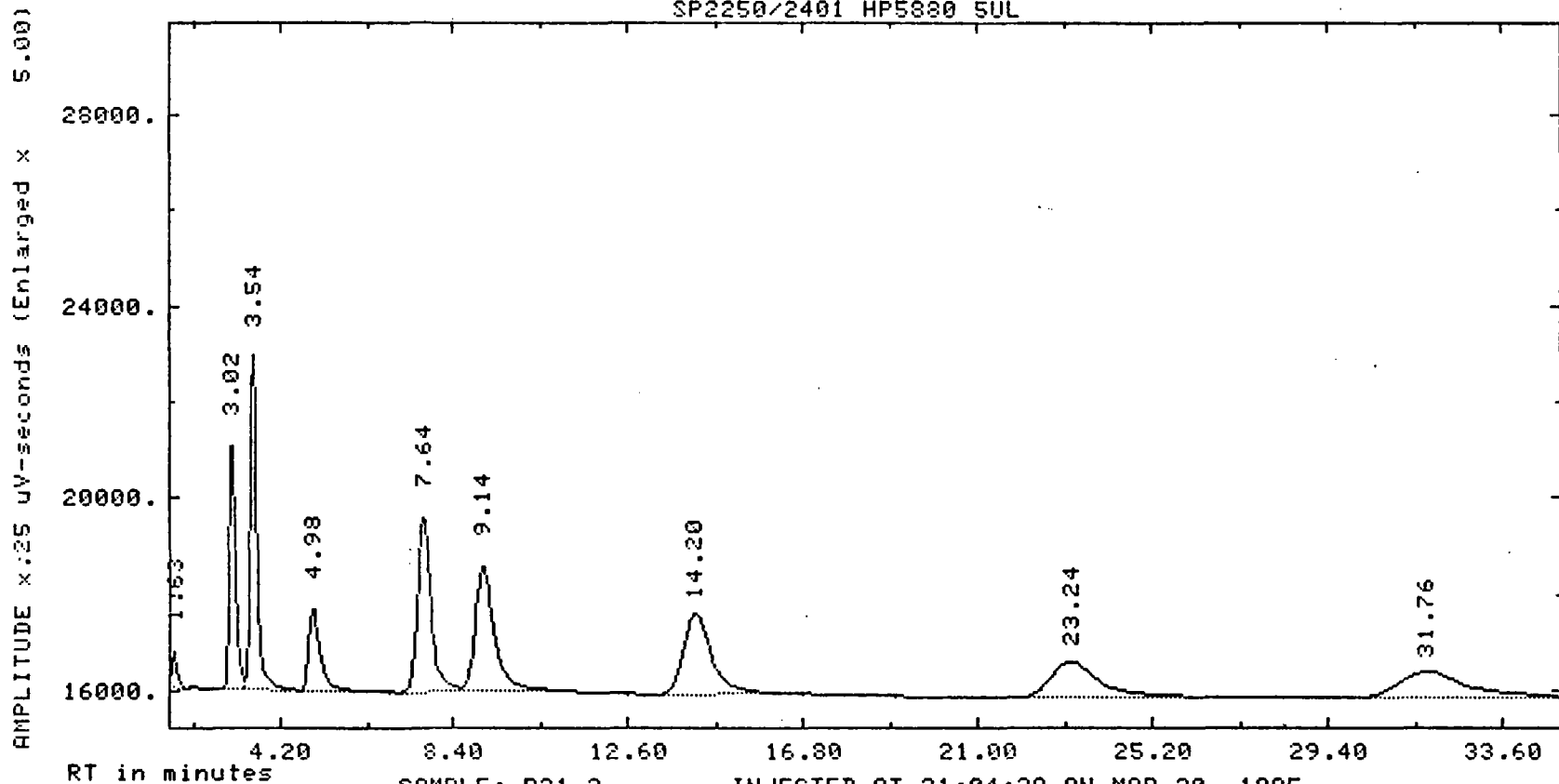
RAW DATA FILE: R5038 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 120

Case 3969

SP2250/2401 HP5880 5UL



AMPLITUDE x:25 uV-seconds (Enlarged x 5.00)

RT in minutes

SAMPLE: P31-3

INJECTED AT 21:04:29 ON MAR 30, 1985

Method: PRIME Raw: R5039 Proc: P5039

121 VI

*LI,P,PS039

PROCESSED DATA FILE: PS039 ON CRN 21

Case 3969

APR 2, 1985 11:45:07

REPORT: 135 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HP5890 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.19	0.00	1.00000	183 BB	.068	
2	1.63	0.00	1.00000	3207 BB	1.195	
3	3.02	0.00	1.00000	26840 BV	10.001	β -BHC
4	3.54	0.00	1.00000	38350 VB	14.514	δ -BHC
5	4.98	0.00	1.00000	16417 BB	6.117	
6	7.54	0.00	1.00000	44101 BV	16.433	Endosulfan I
7	9.14	0.00	1.00000	41152 VB	15.334	pp'-DDE
8	14.20	0.00	1.00000	38997 BB	14.531	Endosulfan II
9	23.24	0.00	1.00000	28871 BB	10.758	Endosulfan Sulfate
10	31.76	0.00	1.00000	29653 BF	11.049	Dibutylchloroacetate

TOTAL AREA = 268371 TOTAL AREA % = 100.000

SAMPLE: P31-3 INJECTED AT 21:04:29 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 39 BTL: 39

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: PS039 PARAM FILES= METHOD: SEQ:

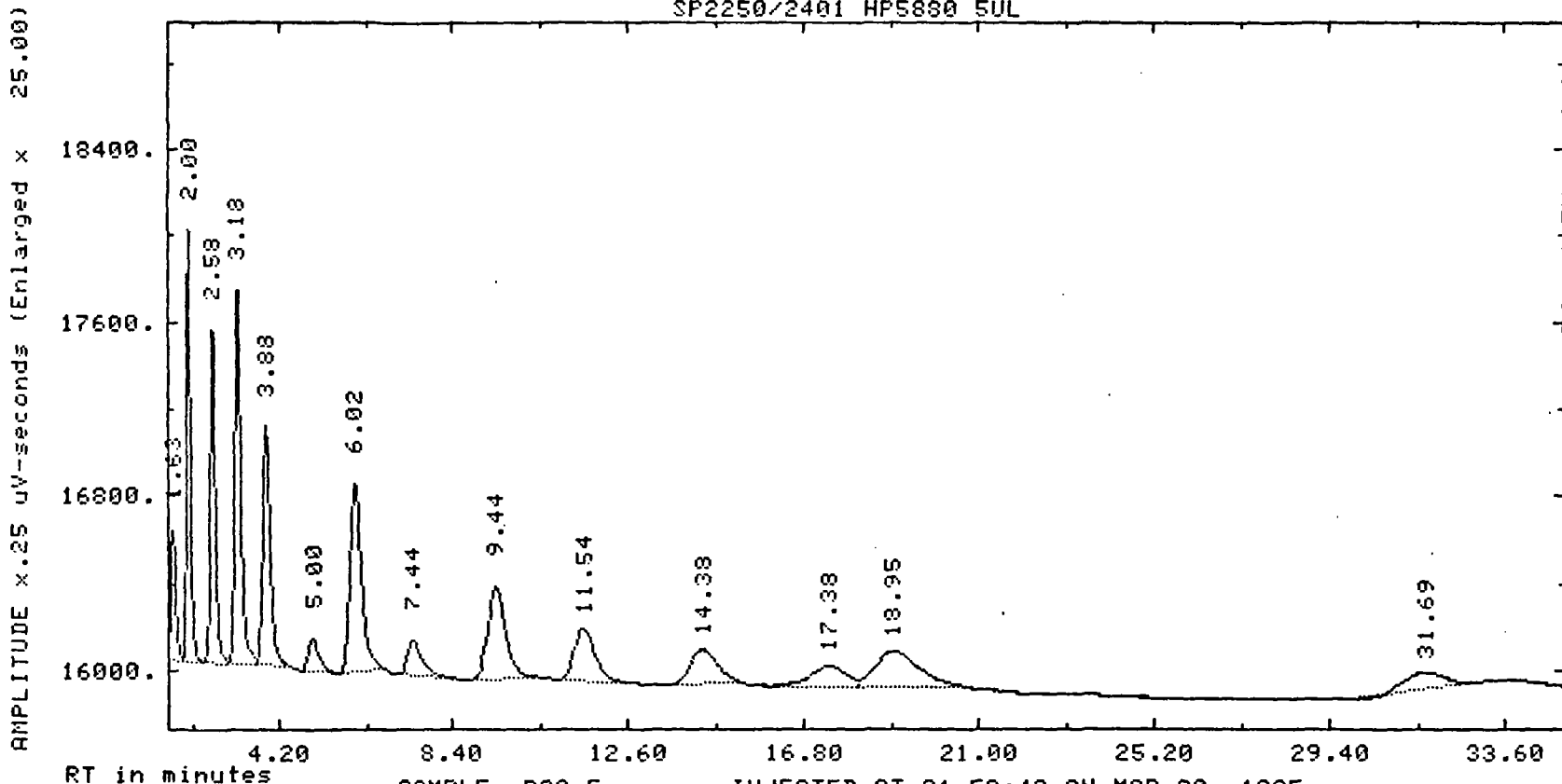
DONE
READY

IV 122

Case 3969

SP2250/2401 HP5880 5UL

301 VI



SAMPLE: P30-5 INJECTED AT 21:59:42 ON MAR 30, 1985
Method: PRIME Raw: R5040 Proc: P5040

*LI,P,P5040

Case 3969

PROCESSED DATA FILE: P5040 ON CRN 21

APR 2, 1985 11:45:31

REPORT: 136 CHANNEL: 2 # PEAKS: -15 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.17	0.00	1.00000	138 BB	.203	
2	1.63	0.00	1.00000	2466 BV	3.640	
3	2.00	0.00	1.00000	6395 VB	9.442	<i>α-BHC</i>
4	2.58	0.00	1.00000	6430 BV	9.484	<i>γ-BHC</i>
5	3.18	0.00	1.00000	9219 VV	13.611	<i>Heptachlor</i>
6	3.88	0.00	1.00000	6961 VB	10.277	<i>Aldrin</i>
7	5.00	0.00	1.00000	1236 BB	1.825	
8	6.02	0.00	1.00000	8180 BB	12.077	<i>Heptachlor Epoxide</i>
9	7.44	0.00	1.00000	2037 BB	3.007	
10	9.44	0.00	1.00000	6132 BB	9.054	<i>Dieldrin</i>
11	11.54	0.00	1.00000	4055 BB	5.996	<i>Endrin</i>
12	14.38	0.00	1.00000	3592 BB	5.304	<i>PP'-DDD</i>
13	17.38	0.00	1.00000	2723 BV	4.021	<i>PP'-DDT</i>
14	18.95	0.00	1.00000	5728 VB	8.457	<i>Endrin Aldehyde</i>
15	31.69	0.00	1.00000	2440 BB	3.602	<i>Endrin Ketone</i>

TOTAL AREA = 67732 TOTAL AREA % = 100.000

SAMPLE: P30-5 INJECTED AT 21:59:42 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 40 BTL: 40

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	XRTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

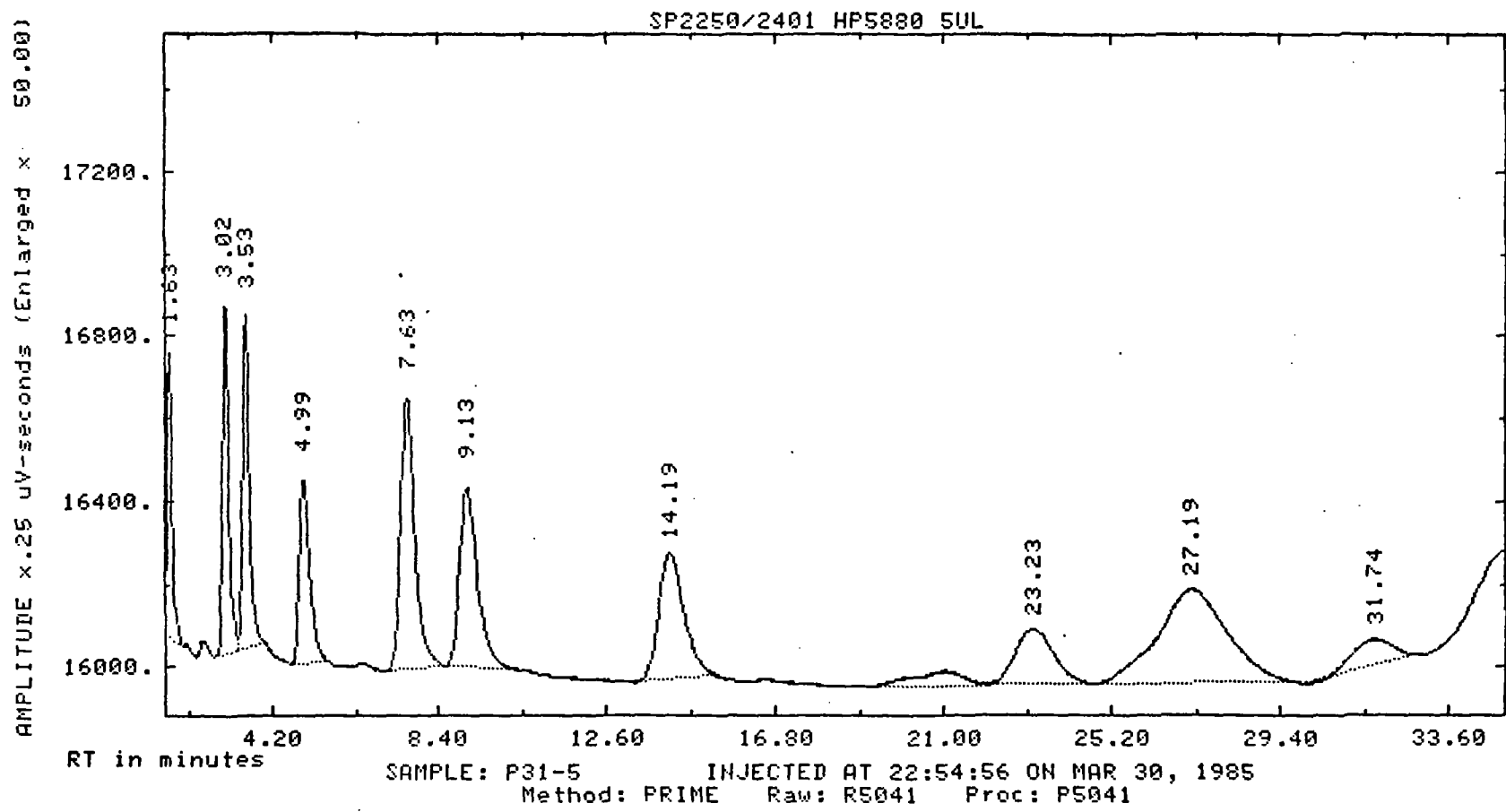
RAW DATA FILE: R5040 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 124

Case 3969

IV
125



*LI,P,P5041

PROCESSED DATA FILE: P5041 ON CRN 21

Case 3969

APR 2, 1985 11:45:58

REPORT: 137 CHANNEL: 2 # PEAKS: -11 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.17	0.00	1.00000	147 BB	.247	
2	1.63	0.00	1.00000	2932 BB	4.936	
3	3.02	0.00	1.00000	4438 BV	7.470	B-BHC
4	3.53	0.00	1.00000	4760 VB	8.012	S-BHC
5	4.99	0.00	1.00000	4008 BB	6.747	
6	7.63	0.00	1.00000	8465 BB	14.249	Endosulfan I
7	9.13	0.00	1.00000	6669 BB	11.226	pp'-DDE
8	14.19	0.00	1.00000	6601 BB	11.214	Endosulfan II
9	23.23	0.00	1.00000	5750 BB	9.679	Endosulfan Sulfate
10	27.19	0.00	1.00000	13443 BB	22.630	
11	31.74	0.00	1.00000	2133 BB	3.591	Dibutylchlorodate

TOTAL AREA = 59405 TOTAL AREA % = 100.000

SAMPLE: P31-5 INJECTED AT 22:54:56 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SMO2 SUBSQ/SAMP: 1 / 41 BTL: 41

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 48.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

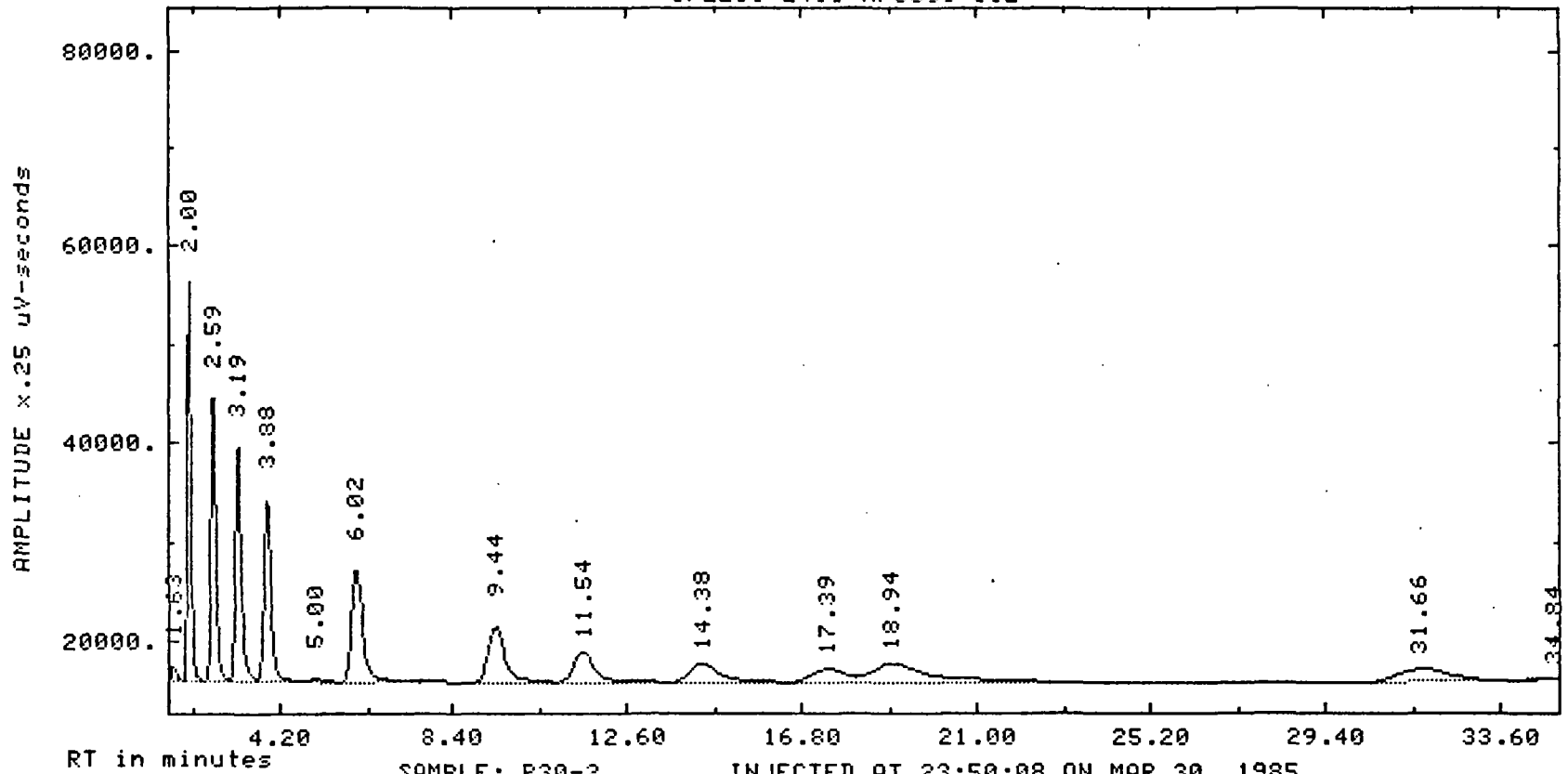
RAW DATA FILE: R5041 PARAM FILES= METHOD: SEQ:

DONE
READY

JV 126

Case 3969

SP2250/2401 HP5890 5UL



SAMPLE: P30-2 INJECTED AT 23:50:08 ON MAR 30, 1985
Method: PRIME Raw: R5042 Proc: P5042

421. III

*LI,P,P5042

PROCESSED DATA FILE: P5042 ON CRN 21

Case 3969

APR 2, 1985 11:46:24

REPORT: 138 CHANNEL: 2 # PEAKS: -17 SP2250/2401 HP5680 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.04	0.00	1.00000	71 BV	.007	
2	1.16	0.00	1.00000	195 VB	.019	
3	1.63	0.00	1.00000	7058 BV	.695	
4	2.00	0.00	1.00000	135844 VV	13.372	<i>α-BHC</i>
5	2.59	0.00	1.00000	120828 VV	11.894	<i>γ-BHC</i>
6	3.19	0.00	1.00000	122525 VV	12.061	<i>Heptachlor</i>
7	3.88	0.00	1.00000	108283 VB	10.659	<i>Aldrin</i>
8	5.00	0.00	1.00000	2903 BB	.286	
9	6.02	0.00	1.00000	105975 BB	10.432	<i>Heptachlor Epoxide</i>
10	9.44	0.00	1.00000	79763 BV	7.852	<i>Dieldrin</i>
11	11.54	0.00	1.00000	54547 VB	5.370	<i>Endrin</i>
12	14.38	0.00	1.00000	47458 BV	4.672	<i>pp'-DDD</i>
13	17.39	0.00	1.00000	37455 VV	3.687	<i>pp'-DDT</i>
14	18.94	0.00	1.00000	87914 VB	8.654	<i>Endrin Aldehyde</i>
15	31.66	0.00	1.00000	54205 BV	5.336	<i>Endrin Ketone</i>
16	34.84	0.00	1.00000	2142 VB	.211	<i>Methoxychlor</i>
17	39.30	0.00	1.00000	48705 BF	4.794	

TOTAL AREA = 1015870 TOTAL AREA % = 100.000

SAMPLE: P30-2 INJECTED AT 23:50:03 ON MAR 30, 1985

ZERO METHOD: PRIME SEQ: SNO3 SUBSQ/SAMP: 1 / 42 BTL: 42

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	XRTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

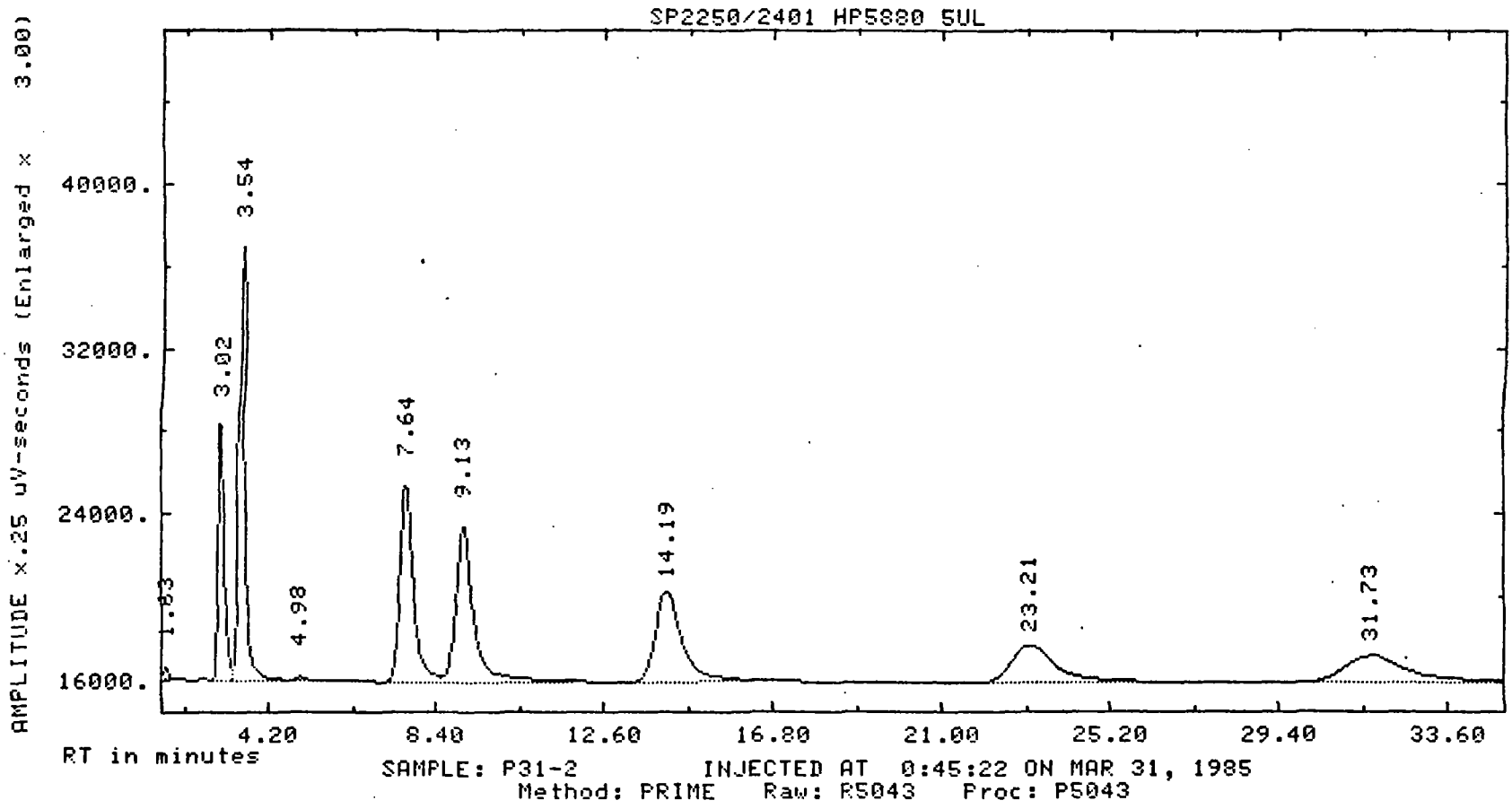
RAW DATA FILE: R5042 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 128

Case 3969

601 AI



*LI,P,P5043

PROCESSED DATA FILE: P5043 ON CRN 21

Case 3969

APR 2, 1985 11:46:55

REPORT: 139 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.20	0.00	1.00000	489 BB	.075	
2	1.63	0.00	1.00000	2349 BB	.359	
3	3.02	0.00	1.00000	66227 BV	10.121	B-BHC
4	3.54	0.00	1.00000	118794 VB	18.154	F-BHC
5	4.98	0.00	1.00000	1605 BD	.245	
6	7.64	0.00	1.00000	111300 BV	17.008	Endosulfan I
7	9.13	0.00	1.00000	115391 VB	17.654	PP-OPE
8	14.19	0.00	1.00000	108621 BB	15.377	Endosulfan II
9	23.21	0.00	1.00000	70727 BB	10.809	
10	31.73	0.00	1.00000	66880 BF	10.220	Dib.tylchloroendate

TOTAL AREA = 654383 TOTAL AREA % = 100.000

SAMPLE: P31-2 INJECTED AT 0:45:22 ON MAR 31, 1985

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 43 BTL: 43

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

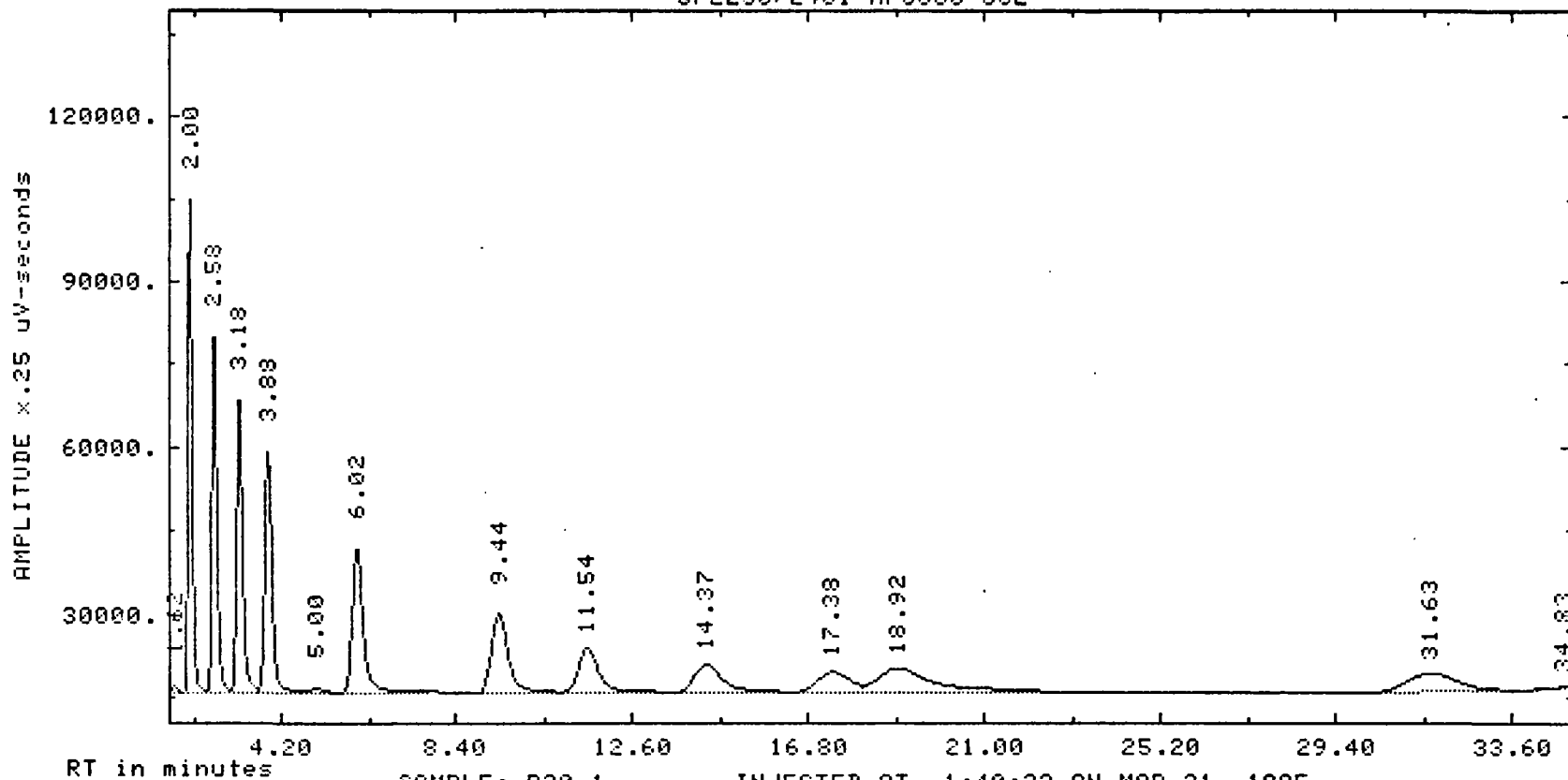
RAW DATA FILE: R5043 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 130

Case 3969

SP2250/2401 HP5880 5UL



RT in minutes

SAMPLE: P30-1

INJECTED AT 1:40:33 ON MAR 31, 1985

Method: PRIME Raw: R5044 Proc: P5044

131

131

*LI,P,P5044

PROCESSED DATA FILE: P5044 ON CRN 21

Case 3969

APR 2, 1985 11:47:19

REPORT: 140 CHANNEL: 2 # PEAKS: -17 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.04	0.00	1.00000	62 BV	.003	
2	1.16	0.00	1.00000	135 VB	.006	
3	1.37	0.00	1.00000	119 BV	.005	
4	1.62	0.00	1.00000	4539 VV	.196	
5	2.00	0.00	1.00000	314054 VV	13.567	α -BHC
6	2.58	0.00	1.00000	263024 VV	12.226	γ -BHC
7	3.18	0.00	1.00000	279720 VV	12.084	Heptachlor
8	3.80	0.00	1.00000	264759 VV	11.437	Aldrin
9	5.00	0.00	1.00000	7872 VB	.340	
10	6.02	0.00	1.00000	244993 BB	10.583	Heptachlor Epoxide
11	9.44	0.00	1.00000	199371 BV	8.559	Dieldrin
12	11.54	0.00	1.00000	141876 VV	6.129	Endrin
13	14.37	0.00	1.00000	125551 VV	5.441	pp'-DDD
14	17.32	0.00	1.00000	103885 VV	4.487	pp'-DDT
15	18.92	0.00	1.00000	202057 VB	8.729	Endrin Aldehyde
16	31.62	0.00	1.00000	132011 BV	5.703	Endrin Ketone
17	34.83	0.00	1.00000	11473 VB	.496	Methoxychlor

TOTAL AREA = 2314083 TOTAL AREA % = 100.000

SAMPLE: P30-1 INJECTED AT 1:40:33 ON MAR 31, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 44 BTL: 44

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

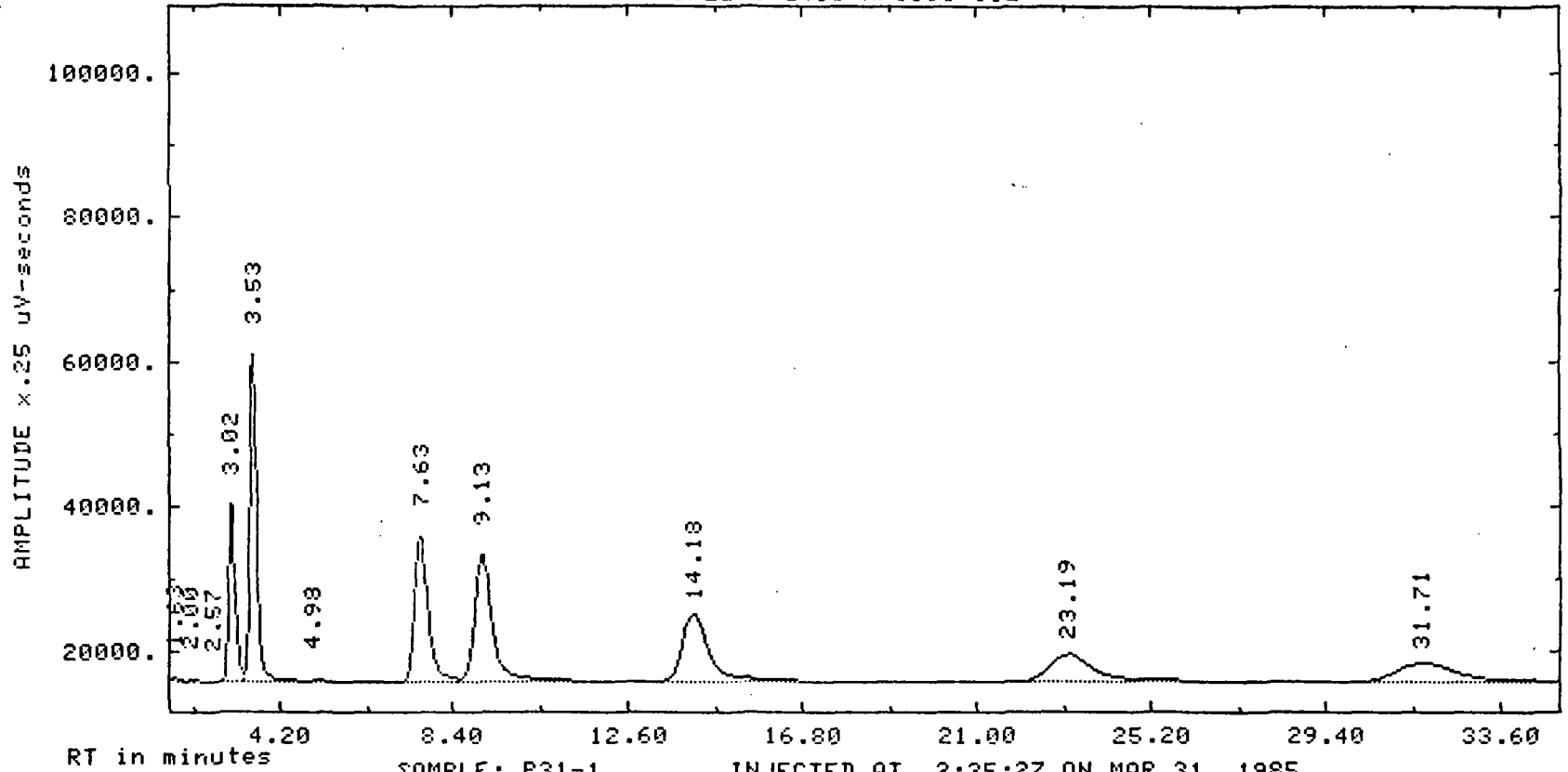
RAW DATA FILE: R5044 PARAM FILES= METHOD: SEQ:

DONE
READY

TV 132

Case 3969

SP2250/2401 HP5880 5UL



SAMPLE: P31-1 INJECTED AT 2:35:27 ON MAR 31, 1985
Method: PRIME Raw: R5045 Proc: P5045

IV 133

*LI,P,P5045

PROCESSED DATA FILE: P5045 ON CRN 21

Case 3969

APR 2, 1985 11:47:48

REPORT: 141 CHANNEL: 2 # PEAKS: -13 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.03	0.00	1.00000	53 BB	.004	
2	1.20	0.00	1.00000	760 BB	.054	
3	1.63	0.00	1.00000	2380 BV	.168	
4	2.00	0.00	1.00000	248 VB	.018	
5	2.57	0.00	1.00000	426 BB	.030	
6	3.02	0.00	1.00000	133339 BV	9.518	B-BHC
7	3.53	0.00	1.00000	257672 VV	18.107	δ-BHC
8	4.93	0.00	1.00000	1715 VB	.122	
9	7.63	0.00	1.00000	235858 BV	16.838	Endosulfan I
10	9.13	0.00	1.00000	264434 VB	18.978	pp'-DDE
11	14.18	0.00	1.00000	215159 BB	15.358	Endosulfan II
12	23.15	0.00	1.00000	148315 BB	10.445	Endosulfan Sulfate
13	31.71	0.00	1.00000	132536 BF	9.461	D. butyl chloroendate

TOTAL AREA = 1400675 TOTAL AREA % = 100.000

SAMPLE: P31-1 INJECTED AT 2:35:27 ON MAR 31, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 45 BTL: 45

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

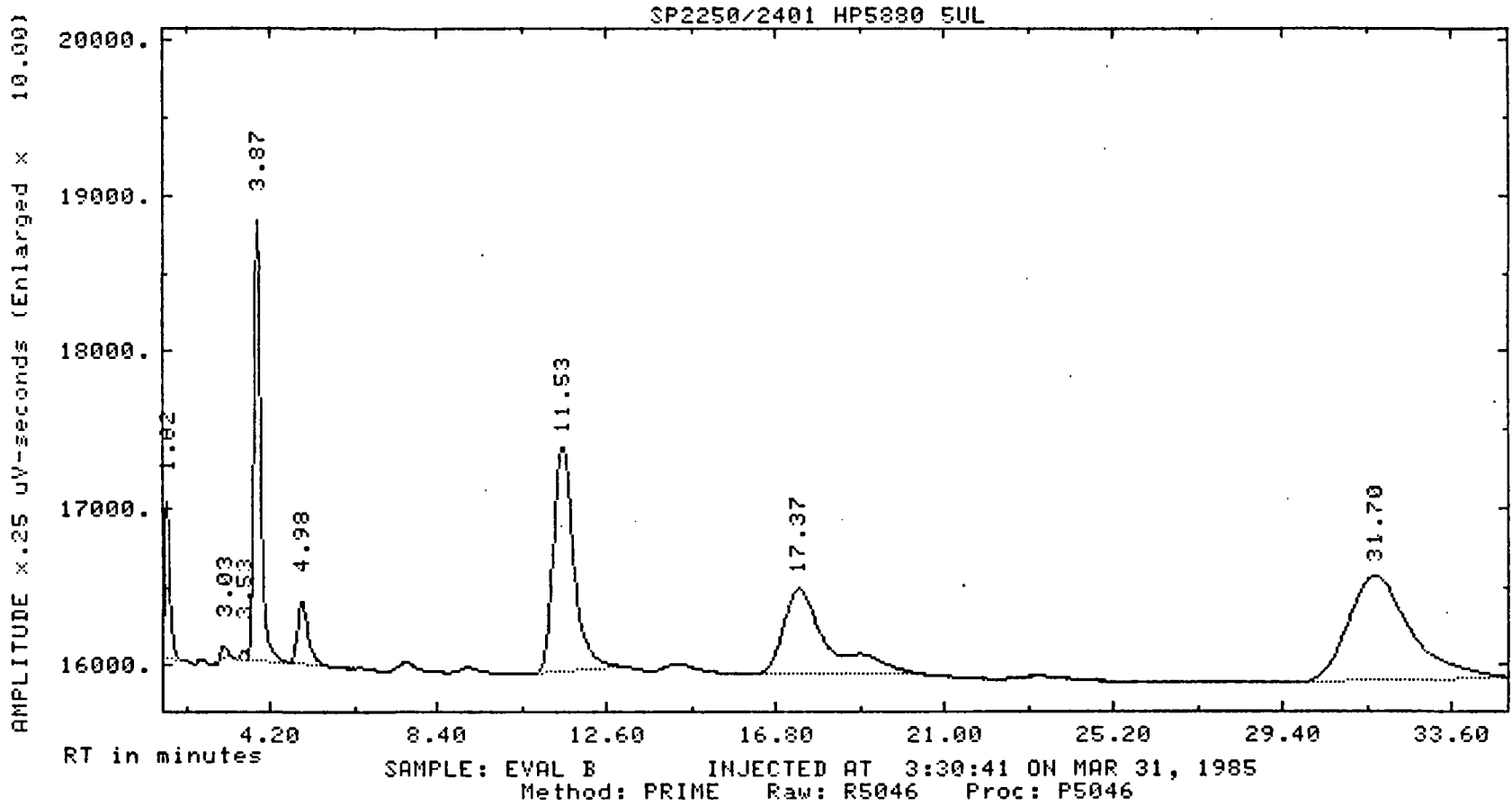
RAW DATA FILE: R5045 PARAM FILES= METHOD: SEQ:

DONE
READY

JW 134

Case 3969

135



*LI,P,P5046

PROCESSED DATA FILE: P5046 ON CRN 21

APR 2, 1985 11:48:11

Case 3969

REPORT: 142 CHANNEL: 2 # PEAKS: -9 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.17	0.00	1.00000	151 BB	.141	
2	1.62	0.00	1.00000	4341 BB	4.059	
3	3.03	0.00	1.00000	391 BB	.365	
4	3.53	0.00	1.00000	284 BV	.266	
5	3.97	0.00	1.00000	16960 VB	15.856	Aldrin
6	4.98	0.00	1.00000	3631 BB	3.394	
7	11.53	0.00	1.00000	24916 BB	23.294	Endrin
8	17.37	0.00	1.00000	20216 BB	18.900	pp'-DDT
9	31.70	0.00	1.00000	36073 BF	33.725	Dibutylchlorodate

TOTAL AREA = 105964 TOTAL AREA % = 100.000

SAMPLE: EVAL B INJECTED AT 3:30:41 ON MAR 31, 1985

ZERO METHOD: PRIME SEQ: SM03 SUBSQ/SAMP: 1 / 46 BTL: 46

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

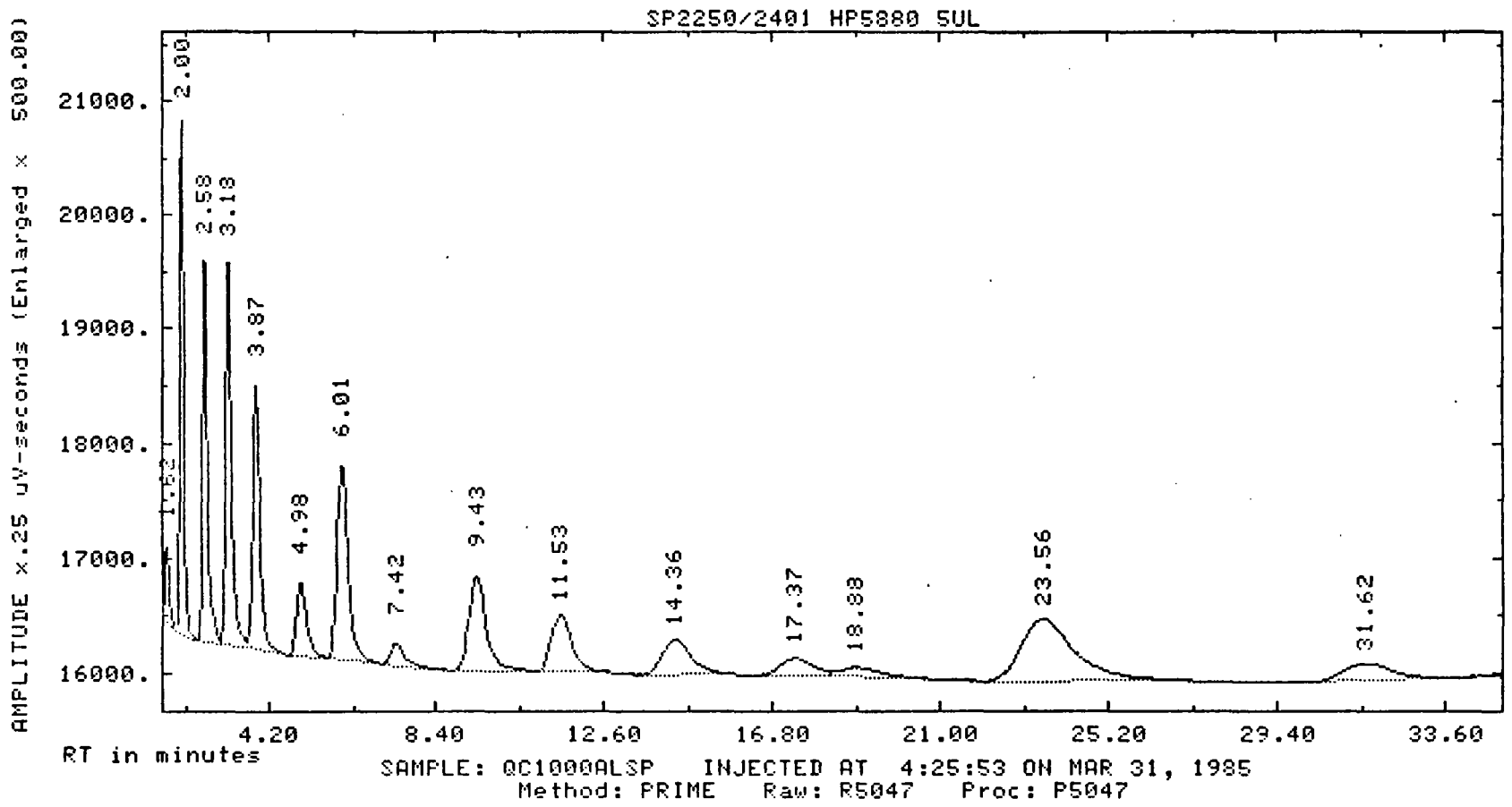
RAW DATA FILE: R5046 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 136

Case 3969

IV
139



*LI,P,P5047

Case 3969

PROCESSED DATA FILE: P5047 ON CRN 21

APR 2, 1985 11:48:34

REPORT: 143 CHANNEL: 2 # PEAKS: -17 SP2250/2401 HPS880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.16	0.00	1.00000	108 BB	.071	
2	1.45	0.00	1.00000	56 BB	.037	
3	1.62	0.00	1.00000	2453 BV	1.612	
4	2.00	0.00	1.00000	14680 VB	9.646	
5	2.58	0.00	1.00000	14127 BV	9.283	
6	3.18	0.00	1.00000	17487 VV	11.491	
7	3.87	0.00	1.00000	14199 VB	9.271	
8	4.98	0.00	1.00000	5909 BB	3.883	
9	6.01	0.00	1.00000	15895 BB	10.510	
10	7.42	0.00	1.00000	2171 BB	1.427	
11	9.43	0.00	1.00000	12023 BB	7.901	
12	11.53	0.00	1.00000	8626 BB	5.668	
13	14.36	0.00	1.00000	7357 BB	4.834	
14	17.37	0.00	1.00000	4304 BV	2.828	
15	18.88	0.00	1.00000	2339 VB	1.537	
16	23.56	0.00	1.00000	24252 BB	15.936	
17	31.62	0.00	1.00000	6187 BB	4.065	

TOTAL AREA = 152185 TOTAL AREA % = 100.000

SAMPLE: QC1000ALSP INJECTED AT 4:25:53 ON MAR 31, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSQ/SAMP: 1 / 47 BTL: 47

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

UNDED NOT ON BASELINE

RAW DATA FILE: R5047 PARAM FILES= METHOD: SEQ:

DONE

IV 138

157

PRIORITY POLLUTANT PESTICIDE DATA

RUN DATE 4/1/85

ANALYST: BL

HP 5880 LOGBOOK PAGE _____

(see reverse for linear regression analyses)

Component	RT (Min)	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std Conc (ng/ml)	Area ₃ X 10 ³	Std Conc (ng/ml)	Area ₃ X 10 ³	Std Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³
1. α-BHC	199-201	100	246.2 ^{2.01}	50	145.4 ^{2.00}	25	58.31 ^{2.00}	10	23.82 ^{2.00}	5	6.96 ^{2.00}		
2. γ-BHC	257-259		270.2 ^{2.59}	(51)	130.2 ^{2.58}	(22.3)	52.12 ^{2.58}	(11.1)	21.75 ^{2.58}	(5.5)	6.81 ^{2.58}		
3. Heptachlor	317-319	(97.1)	227.9 ^{3.19}	(56.6)	134.6 ^{3.18}	(23.3)	52.30 ^{3.18}	(10.4)	22.14 ^{3.18}	(2.3)	9.41 ^{3.18}		
4. Aldrin	386-370		217.4 ^{3.71}	(54.0)	117.6 ^{3.78}	(22.1)	46.49 ^{3.78}	(10.8)	21.14 ^{3.78}	(4.4)	6.87 ^{3.78}		
5. Hept. Epox.	571-605		187.6 ^{6.03}		113.6 ^{6.01}		48.93 ^{6.01}		24.10 ^{6.02}		10.04 ^{6.01}		
6. Dieldrin	741-745	(99.0)	162.3 ^{7.44}	(53.1)	87.26 ^{7.43}	(22.5)	37.07 ^{7.42}	(10.5)	17.46 ^{7.43}	(4.9)	8.28 ^{7.43}		
7. Endrin	1149-1155	(100.4)	118.7 ^{11.54}	(50.6)	59.78 ^{11.52}	(21.6)	25.48 ^{11.52}	(10.6)	12.44 ^{11.53}	(6.8)	8.02 ^{11.51}		
8. p,p'-DDD	1428-1440		77.0 ^{14.31}		50.44 ^{14.34}		20.19 ^{14.34}		10.30 ^{14.35}		12.64 ^{14.31}		
9. p,p'-DDT	1729-1741	(86.5)	67.89 ^{17.38}	(57.4)	44.72 ^{17.34}	(21.1)	18.65 ^{17.33}	(8.0)	7.73 ^{17.35}		-		
10. Endrin Ald.	1883-1895		140.7 ^{18.91}		94.17 ^{18.87}		42.42 ^{18.87}		20.12 ^{18.74}				
11. Endrin Ketone	3133-3183		121 ^{31.69}		79.91 ^{31.57}		34.77 ^{31.56}		17.91 ^{31.60}		4.48 ^{31.46}		
12. Methoxychlor	3452-3504	↓	60.13 ^{34.90}	↓	43.84 ^{34.75}	↓	18.88 ^{34.71}	↓	9.22 ^{34.74}	↓			
Pesticide Mix		P-30-1		P-30-2		P-30-3		P-30-4		P-30-5			

131

LINEAR REGRESSION RESULTS

SP2250/2401

4-1-85

Component	Correlation Coefficient	Y-intercept	Slope
1. γ -BHC	.9960 50-5ppb stds	-9.20	2.71
2. -BHC			
3. Heptachlor	.9949	4.15	2.30
4. Aldrin	.9978	-2.97	2.23
5. Heptachlor Epoxide			
6. Dieldrin	.9985	.302	1.64
7. Endrin	.9987	-0.06	1.18
8. p,p'-DDD			
9. p,p'-DDT	.987 Highest four stds	4.29	0.68
10. Endrin Aldehyde			
11. Endrin Ketone			
12. Methoxychlor			

149

PRIORITY POLLUTANT PESTICIDE DATA

RUN DATE: 4/1/95

ANALYST: BL

HP 5880 LOGBOOK PAGE _____

(see reverse for linear regression analyses)

Component	RT (Min)	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³	Std. Conc (ng/ml)	Area ₃ X 10 ³
1. α -BHC	2.99-3.05	100	3.04 72.81	50 ³	3.02 60.43	25	3.02 27.40	10	3.02 9.20	5	3.02 4.37		
2. β -BHC	3.50-3.58		3.56 177.7		3.53 157.		3.53 40.63		3.53 10.71		3.53 4.82		
3. Endosulfan I	7.56-7.70		7.67 183.2		7.63 104.3	(23.5)	7.63 45.27	(8.1)	7.63 16.03	(4.9)	7.61 10.04		
4. p,p'-DDE	9.04-9.22		9.15 187.1		9.11 101.2		9.11 41.39		9.12 13.90		9.12 6.55		
5. Endosulfan II	14.08-14.28		14.24 152.8		14.16 91.57		14.16 39.72		14.17 14.26		14.16 6.71		
6. Endosulfan Sulfate	23.02-23.34		23.27 327.8		23.14 67.50		23.14 32.74		23.17 11.21		23.16 6.04		
7. Dibutyl Chlorendate	31.53-31.83	↓ (101)	31.77 137.6	↓ (46.9)	31.64 59.70	↓ (26.6)	31.66 32.40	↓ (10.8)	31.68 11.15	↓ (4.6)	31.66 2.85		
Pesticide Mix			P-31-1		P-31-2		P-31-3		P-31-4		P-31-5		

11:11

LINEAR REGRESSION RESULTS

0.2250/2401

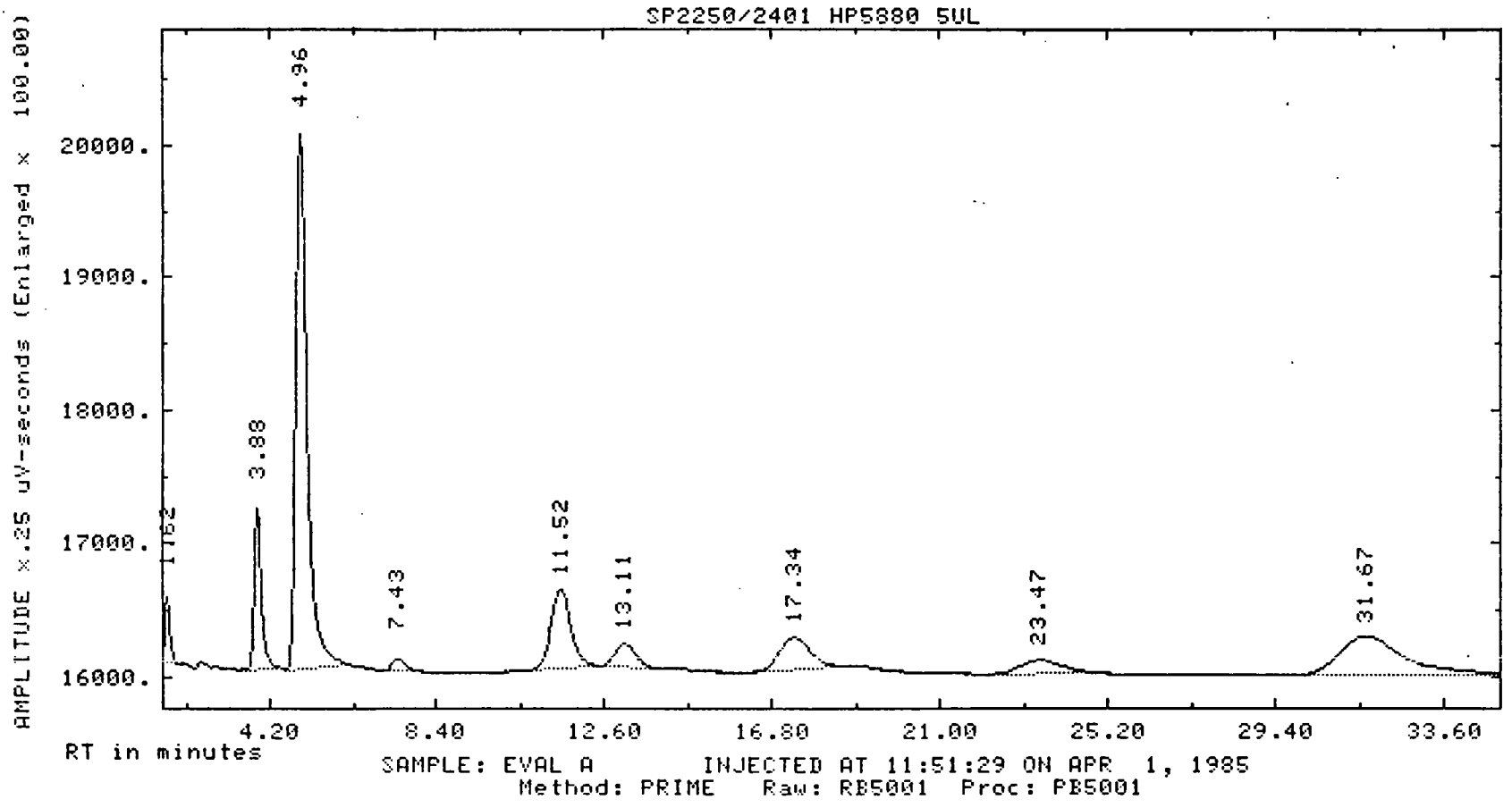
4/1/85

Component	Correlation Coefficient	Y-intercept	Slope
1. δ -BHC			
2. δ -BHC			
3. Endosulfan I			
4. p'-DDE			
5. Endosulfan II			
6. Endosulfan Sulfate			
7. Dibutyl Chlorendate	.9988	-3.36	1.34

IV
142

Case 3969

IV 143



READY
*LI,P,FB5001

PROCESSED DATA FILE: FB5001 ON CRN 21

APR 2, 1985 8:57:15

Case 3969

REPORT: 144 CHANNEL: 2 # PEAKS: -9 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	1975 BB	2.199	
2	3.88	0.00	1.00000	7537 BB	8.368	Aldrin
3	4.96	0.00	1.00000	40232 BB	44.774	
4	7.43	0.00	1.00000	1075 BB	1.197	
5	11.52	0.00	1.00000	10273 BB	11.432	Endrin
6	13.11	0.00	1.00000	3183 BB	3.543	
7	17.34	0.00	1.00000	6241 BB	6.945	pp'-DDT
8	23.47	0.00	1.00000	3421 BB	3.874	
9	31.67	0.00	1.00000	15858 BB	17.643	Dibutylchloroendate

TOTAL AREA = 89856 TOTAL AREA % = 100.000

SAMPLE: EVAL A INJECTED AT 11:51:29 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 1 BTL: 1

SL-WIDTH MV/MIN DELAY MIN-AR BUNCH REPORT
.500 .300 1.00 50 AUTO MEDIUM

SUP-UNK DVT ID-LVL REF-RTW %RTW %DIL-F ISO
NO 0.00 0 .30 5.0 100.00 YES

ACTUAL RUN TIME: 40.008 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: FB5001 PARAM FILES: METHOD: SEQ:

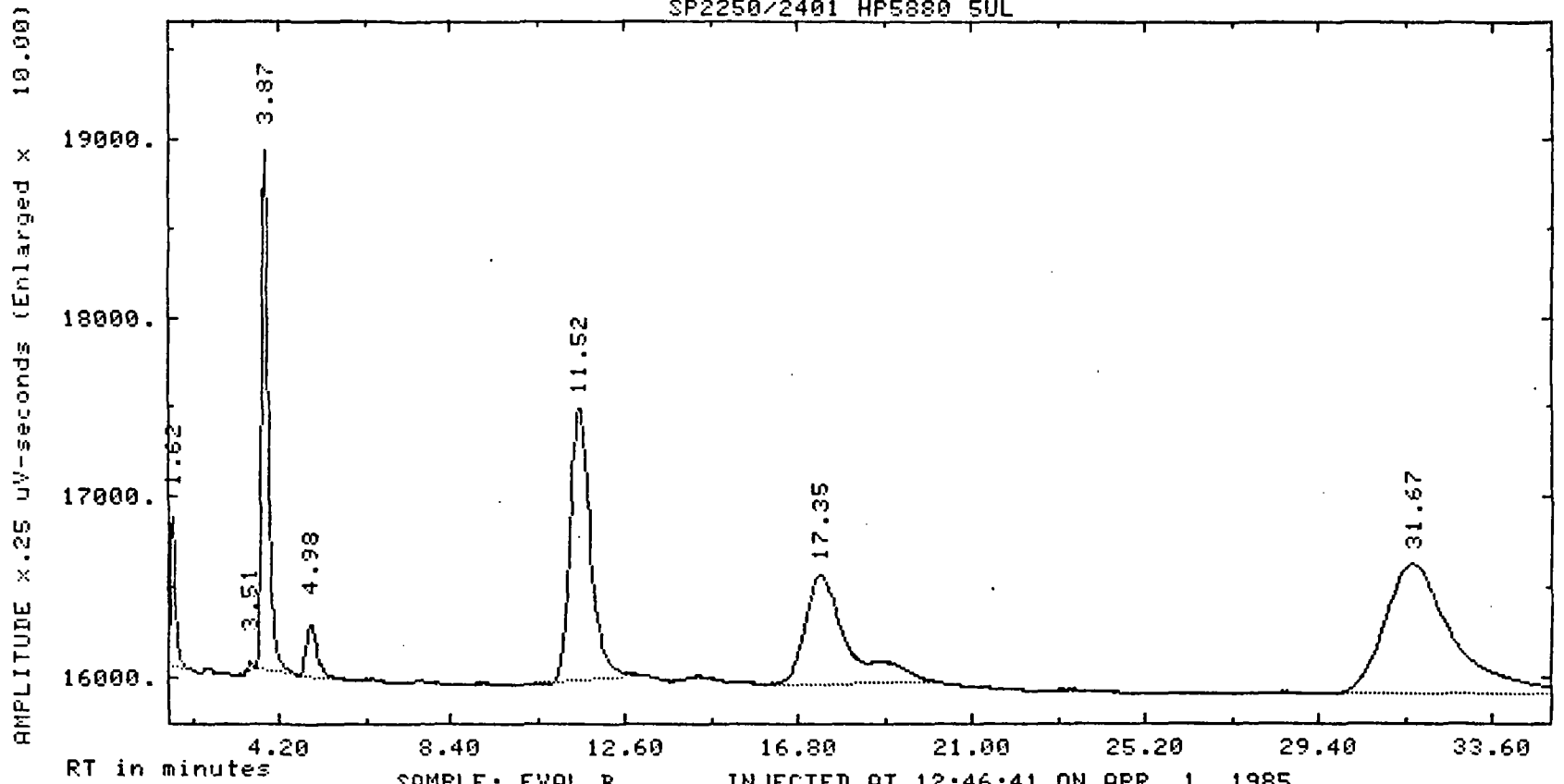
DONE
READY

IV 144

Case 3969

SHI VI

SP2250/2401 HP5880 5UL



AMPLITUDE x.25 uV-seconds (Enlarged x 10.00)

RT in minutes

SAMPLE: EVAL B INJECTED AT 12:46:41 ON APR 1, 1985
Method: PRIME Raw: RB5002 Proc: PB5002

*LI,P,PB5002

PROCESSED DATA FILE: PB5002 ON CRN 21

Case 3969

APR 2, 1985 8:57:39

REPORT: 145 CHANNEL: 2 # PEAKS: -7 SF2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	3469 BB	3.103	
2	3.51	0.00	1.00000	114 BB	.102	
3	3.87	0.00	1.00000	17058 BB	15.256	Aldrin
4	4.98	0.00	1.00000	2558 BB	2.280	
5	11.52	0.00	1.00000	25526 BB	22.833	Endrin
6	17.35	0.00	1.00000	21636 BB	19.353	pp'-DDT
7	31.67	0.00	1.00000	41436 BB	37.064	D.b.tylochloredate

TOTAL AREA = 111757 TOTAL AREA % = 100.000

SAMPLE: EVAL B INJECTED AT 12:46:41 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 2 BTL: 2

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: RE5002 PARAM FILES= METHOD: SEQ:

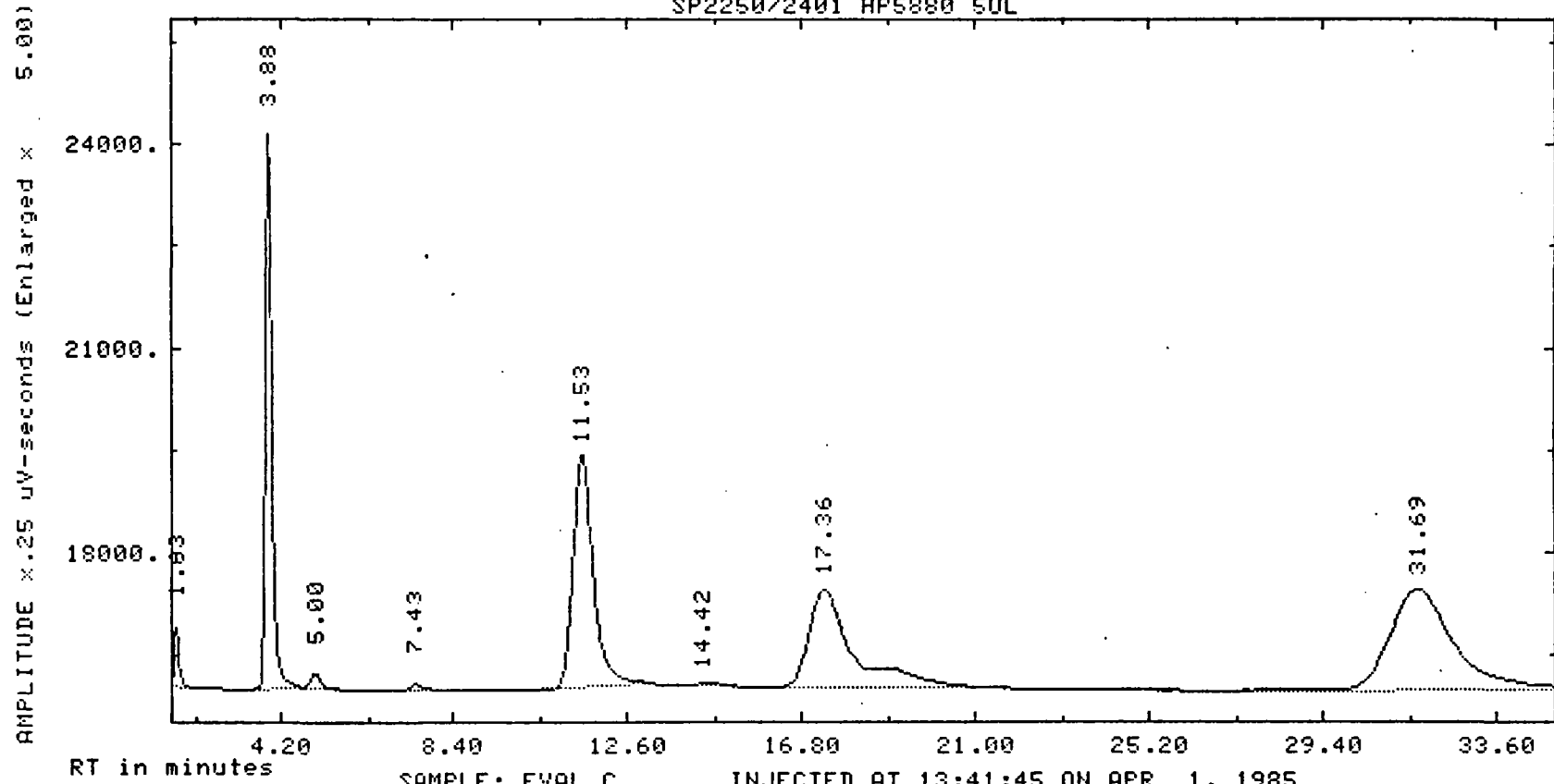
DONE
READY

JV 146

Case 3969

SP2250/2401 HP5880 5UL

LH1 IR



RT in minutes
SAMPLE: EVAL C INJECTED AT 13:41:45 ON APR 1, 1985
Method: PRIME Raw: RB5003 Proc: PB5003

*L1,P,PD5003

PROCESSED DATA FILE: PD5003 ON CRN 21

Case 3969

APR 2, 1985 8:58:02

REPORT: 146 CHANNEL: 2 # PEAKS: -8 SP2250/2401 HP5880 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	3588 BB	1.499	
2	3.88	0.00	1.00000	46327 BB	19.357	Aldrin
3	5.00	0.00	1.00000	1887 BB	.788	
4	7.43	0.00	1.00000	993 BB	.415	
5	11.53	0.00	1.00000	56591 BB	23.645	Endrin
6	14.42	0.00	1.00000	605 BB	.286	
7	17.36	0.00	1.00000	52098 BB	21.786	pp'-DDT
8	31.89	0.00	1.00000	77165 BB	32.242	Dibutylchlorodate

TOTAL AREA = 238335 TOTAL AREA % = 100.000

SAMPLE: EVAL C INJECTED AT 13:41:45 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 3 BYL: 3

SL-WIDTH	MOV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

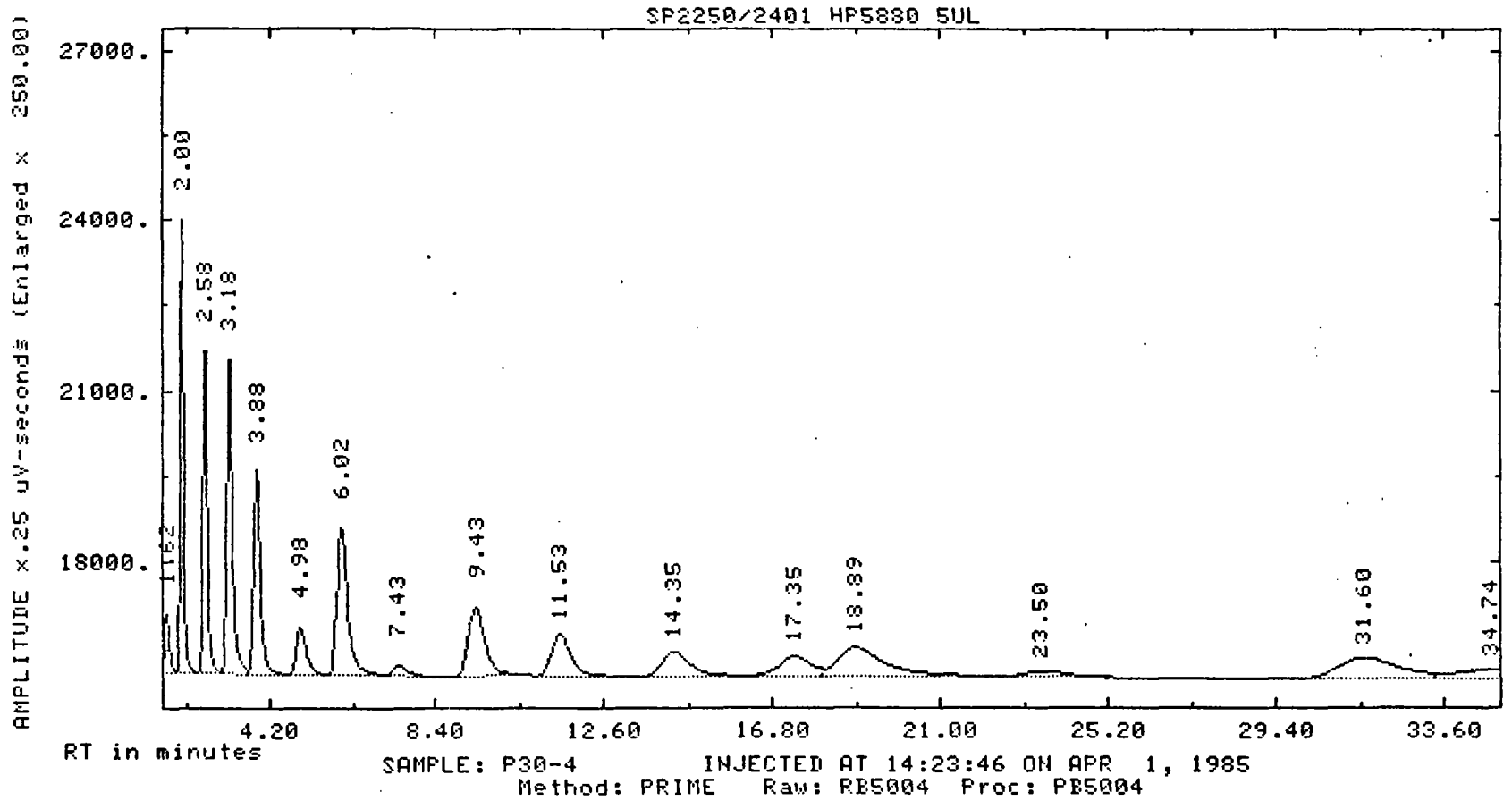
RAW DATA FILE: PD5003 PARAM FILES= METHOD: SEQ:

DONE
READY

JV 148

Case 3969

641 II



*LI,P,PB5004

PROCESSED DATA FILE: PB5004 ON CRN 21

Case 3969

APR 2, 1985 8:58:28

REPORT: 147 CHANNEL: 2 # PEAKS: -16 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	4159 BV	1.783	
2	2.00	0.00	1.00000	23018 VB	10.212	<i>α-BHC</i>
3	2.58	0.00	1.00000	21751 BV	9.326	<i>γ-BHC</i>
4	3.18	0.00	1.00000	20135 VV	12.053	<i>Heptachlor</i>
5	3.88	0.00	1.00000	21135 VB	9.062	<i>Aldrin</i>
6	4.98	0.00	1.00000	7340 BB	3.147	
7	5.02	0.00	1.00000	24097 BB	10.332	<i>Heptachlor Epoxide</i>
8	7.43	0.00	1.00000	2209 BB	.947	
9	9.43	0.00	1.00000	17457 BB	7.485	<i>pp'-DDT BL. Dieldrin</i>
10	11.53	0.00	1.00000	12438 BB	5.332	<i>Endrin</i>
11	14.35	0.00	1.00000	10304 BB	4.418	<i>pp'-DDD</i>
12	17.35	0.00	1.00000	9727 BV	4.171	<i>pp'-DDT</i>
13	18.89	0.00	1.00000	20121 VB	8.627	<i>Endrin Aldehyde</i>
14	23.50	0.00	1.00000	3405 BB	1.460	
15	31.60	0.00	1.00000	17814 BV	7.681	<i>Endrin Ketone</i>
16	34.74	0.00	1.00000	9221 VB	3.954	<i>Methoxychlor</i>

TOTAL AREA = 233232 TOTAL AREA % = 100.000

SAMPLE: P30-4 INJECTED AT 14:23:46 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 4 BTL. 4

SL-WIDTH	NO/MIN	DELAY	MIN-AP	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTN	XRTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

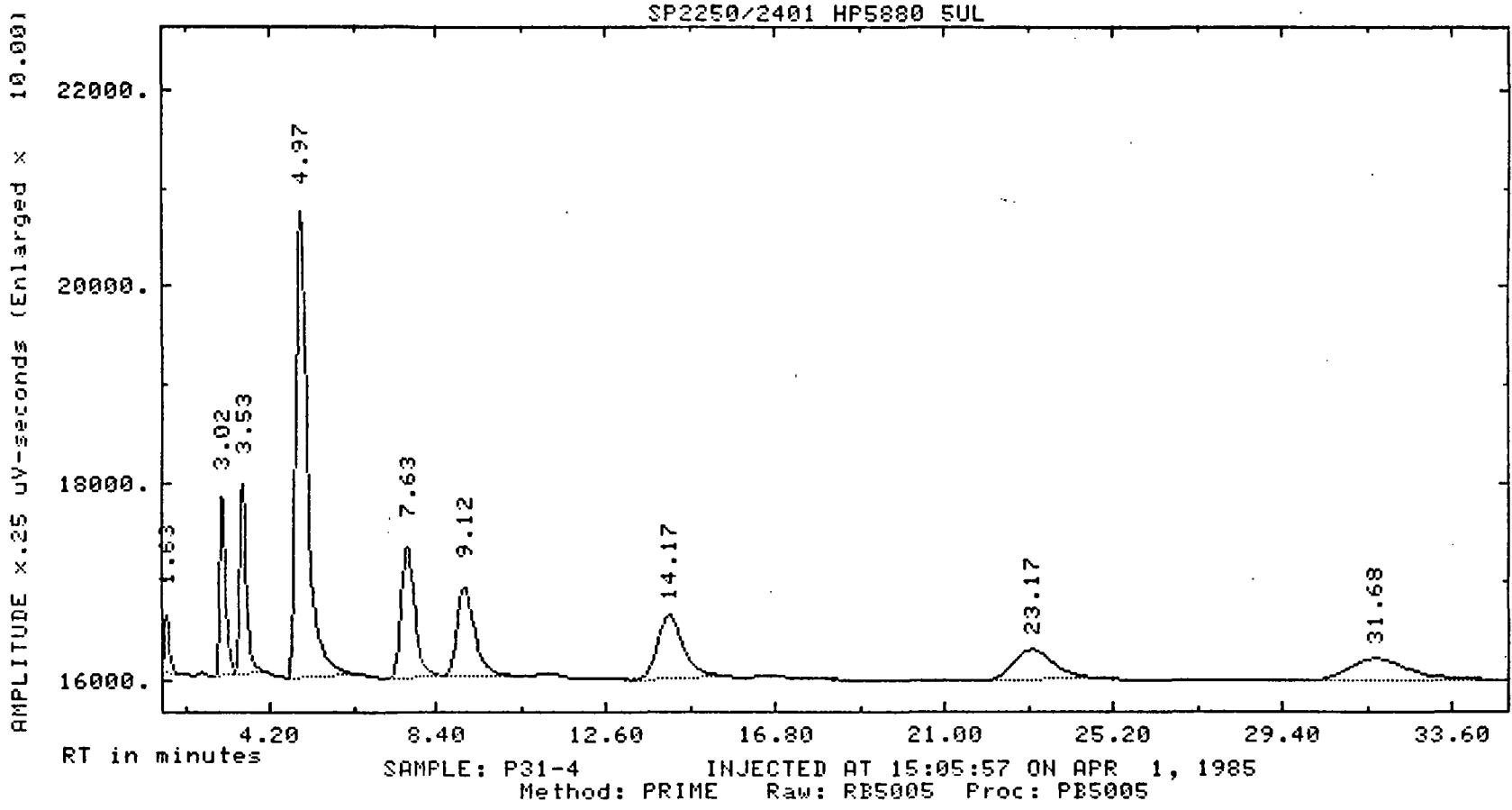
RAW DATA FILE: RE5004 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 150

Case 3969

151 IV



*LI,P,PB5005

PROCESSED DATA FILE: PB5005 ON CRN 21

Case 3969

APR 2, 1985 8:59:00

REPORT: 148 CHANNEL: 2 # PEAKS: -9 SP2250/2401 HP5000 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.63	0.00	1.00000	2451 BB	1.778	
2	3.02	0.00	1.00000	9198 BV	6.671	B-BHC
3	3.53	0.00	1.00000	10785 VB	7.621	σ-BHC
4	4.97	0.00	1.00000	48902 BB	35.465	
5	7.63	0.00	1.00000	16033 BV	11.628	Endosulfan I
6	9.12	0.00	1.00000	13900 VB	10.060	pp'-DDE
7	14.17	0.00	1.00000	14261 BB	10.342	Endosulfan II
8	23.17	0.00	1.00000	11208 BB	8.128	Endosulfan Sulfate
9	31.68	0.00	1.00000	11151 BB	8.087	D:butyl chloroendate

TOTAL AREA = 137890 TOTAL AREA % = 100.000

SAMPLE: P21-4 INJECTED AT 15:05:57 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 5 BTL: 5

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

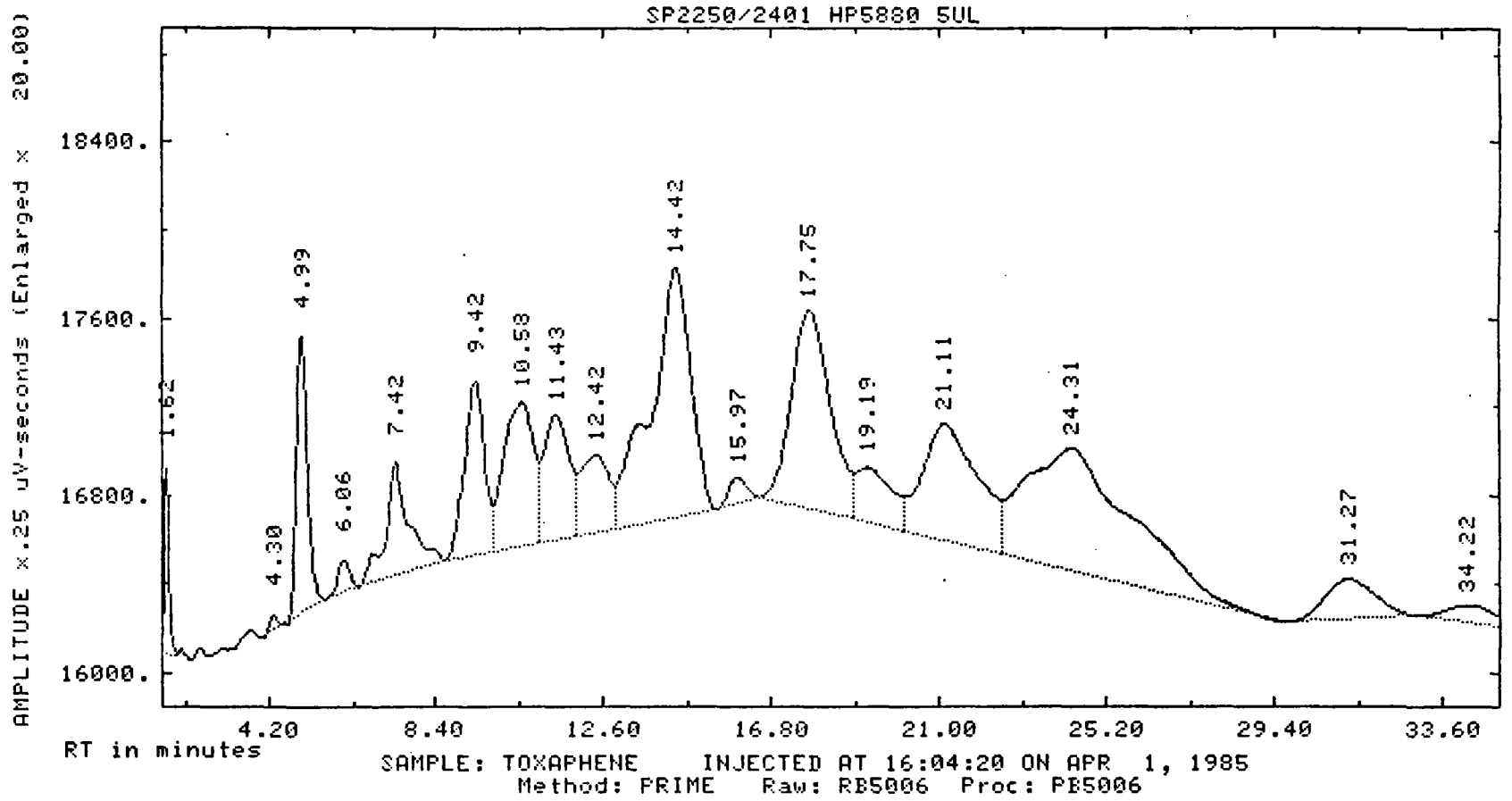
RAW DATA FILE: RB5005 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 152

Case 3969

551 AF



*LI,P,PB5006

PROCESSED DATA FILE: PB5006 ON CRN 21

Case 3969

APR 2, 1985 8:59:30

REPORT: 149 CHANNEL: 2 # PEAKS: -17 SP2250/2401 HPS280 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	3350 BB	1.369	
2	4.30	0.00	1.00000	315 BB	.129	
3	4.99	0.00	1.00000	11984 BB	4.895	
4	6.06	0.00	1.00000	1464 BB	.590	
5	7.42	0.00	1.00000	10444 BB	4.257	
6	9.42	0.00	1.00000	13610 BV	5.560	
7	10.59	0.00	1.00000	16676 VV	6.813	
8	11.43	0.00	1.00000	12506 VV	5.109	
9	12.42	0.00	1.00000	8249 VV	3.370	
10	14.42	0.00	1.00000	39254 VV	16.037	
11	15.97	0.00	1.00000	1814 VB	.741	
12	17.75	0.00	1.00000	30482 BV	12.453	
13	19.19	0.00	1.00000	7764 VV	3.172	
14	21.11	0.00	1.00000	25605 VV	10.451	
15	24.31	0.00	1.00000	50459 VV	20.615	
16	31.27	0.00	1.00000	7689 VB	3.141	
17	34.22	0.00	1.00000	3106 BB	1.269	

TOTAL AREA = 244770 TOTAL AREA % = 100.000

SAMPLE: TOXAPHENE INJECTED AT 16:04:20 ON APR 1, 1985

ZERO METHOD: PRINL SEQ: SMO3A SUBSQ/SAMP: 1 / 6 BTL: 6

SL-WDTH	NU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.025 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

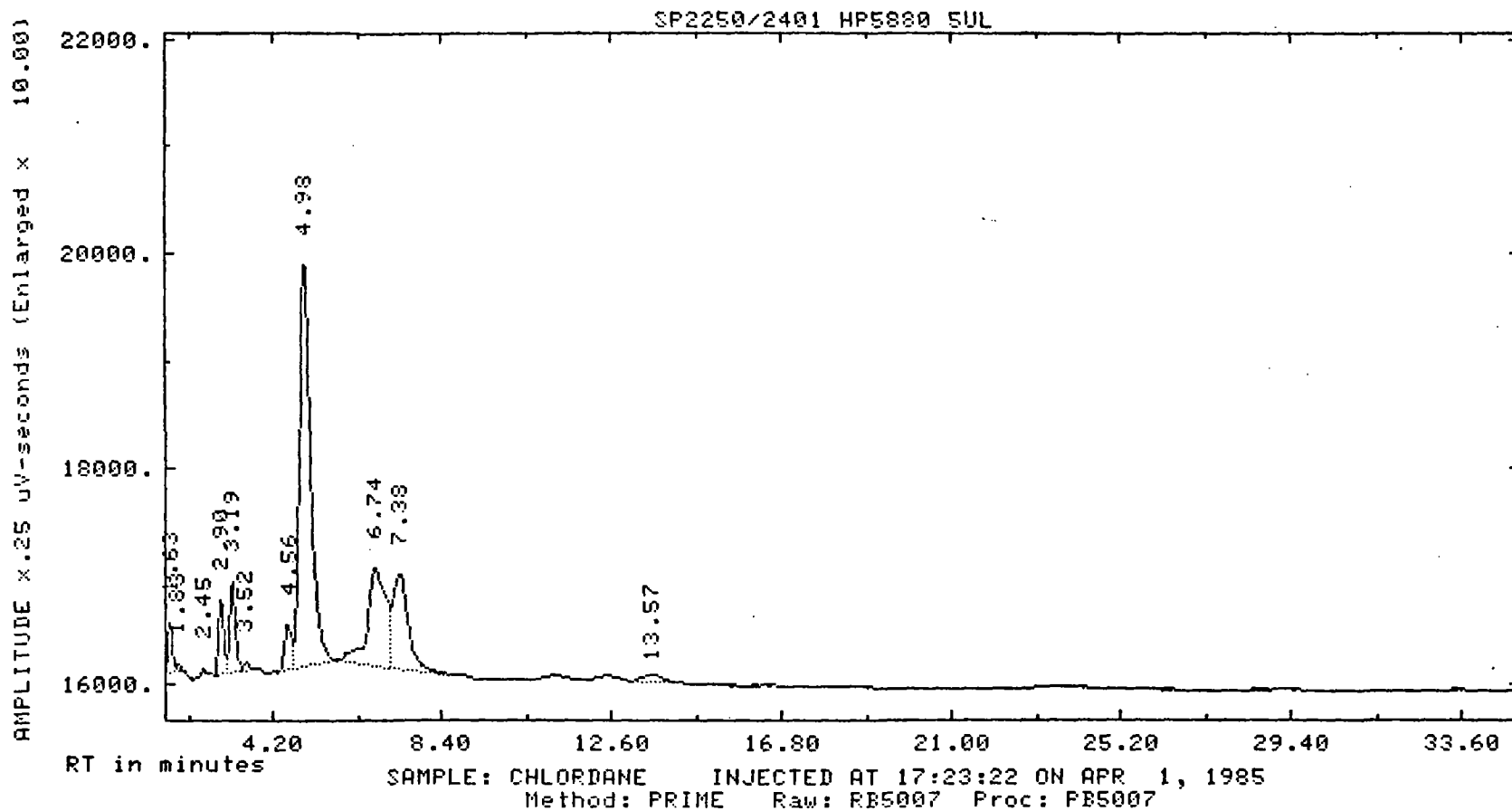
RAW DATA FILE: RB5006 PARAM FILES= METHOD: SEQ:

DONE
READY

TU 154

Case 3969

551 IV



*LI,P,P85007

PROCESSED DATA FILE: P85007 ON CRN 21

Case 3969

APR 2, 1985 8:59:59

REPORT: 150 CHANNEL: 2 # PEAKS: -12 SP2250/2401 HFS880 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.30	0.00	1.00000	658 BB	.862	
2	1.63	0.00	1.00000	1737 BV	2.240	
3	1.83	0.00	1.00000	181 VB	.234	
4	2.45	0.00	1.00000	198 BB	.255	
5	2.90	0.00	1.00000	3079 BV	3.971	
6	3.19	0.00	1.00000	4357 VV	5.619	
7	3.52	0.00	1.00000	390 VB	.503	
8	4.56	0.00	1.00000	2649 BV	3.417	
9	4.98	0.00	1.00000	36779 VB	47.437	
10	6.74	0.00	1.00000	15514 BV	20.009	
11	7.38	0.00	1.00000	10993 VB	14.186	
12	13.57	0.00	1.00000	982 BB	1.268	

TOTAL AREA = 77533 TOTAL AREA % = 100.000

SAMPLE: CHLORDANE INJECTED AT 17:23:22 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMOZA SUBSQ/SAMP: 1 / 7 BTL: 7

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-ANT: 1.0000 SAMP-ANT: 1.0000

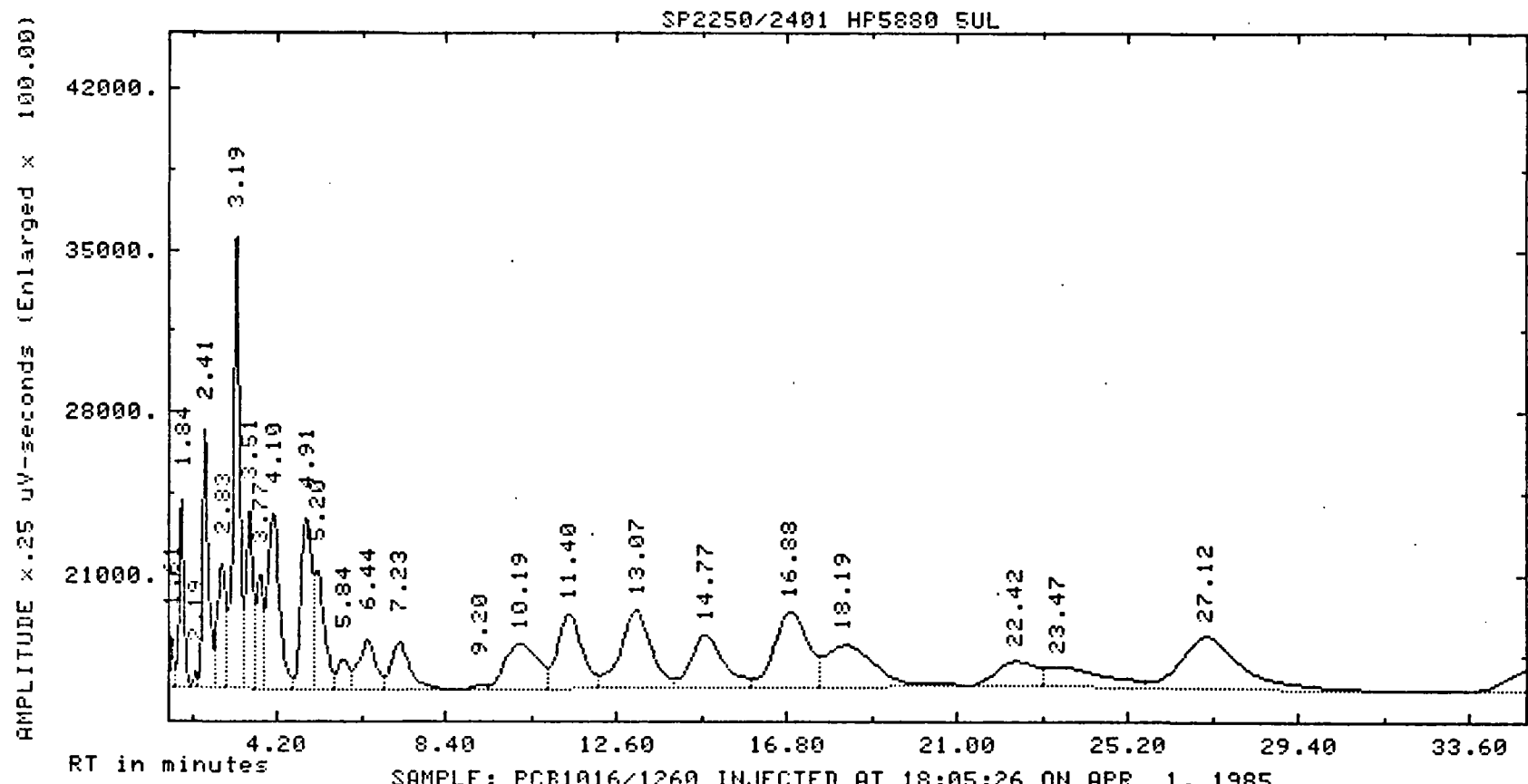
RAW DATA FILE: P85007 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 156

Case 3969

LSI II



SAMPLE: PCB1016/1260 INJECTED AT 18:05:26 ON APR 1, 1985
Method: PRIME Raw: RB5008 Proc: FB5008

ALI, P, P85008

PROCESSED DATA FILE: P85008 ON CRN 21

Case 3969

APR 2, 1985 9:00:23

REPORT: 151 CHANNEL: 2 # PEAKS: -26 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.46	0.00	1.00000	2458 BV	.184	
2	1.61	0.00	1.00000	6991 UV	.523	
3	1.84	0.00	1.00000	32391 UV	2.423	
4	2.19	0.00	1.00000	2588 UV	.194	
5	2.41	0.00	1.00000	56015 UV	4.190	
6	2.63	0.00	1.00000	35324 UV	2.642	
7	3.19	0.00	1.00000	116263 UV	8.696	
8	3.51	0.00	1.00000	47090 UV	3.507	
9	3.77	0.00	1.00000	28199 UV	2.109	
10	4.10	0.00	1.00000	70261 UV	5.854	
11	4.91	0.00	1.00000	70051 UV	5.240	
12	5.20	0.00	1.00000	43339 UV	3.287	
13	5.84	0.00	1.00000	12622 UV	.944	
14	6.44	0.00	1.00000	31012 UV	2.320	
15	7.23	0.00	1.00000	31557 VB	2.360	
16	9.20	0.00	1.00000	2393 BV	.179	
17	10.19	0.00	1.00000	55042 UV	4.117	
18	11.40	0.00	1.00000	67265 UV	5.031	
19	13.07	0.00	1.00000	87557 UV	6.550	
20	14.77	0.00	1.00000	64203 UV	4.802	
21	16.98	0.00	1.00000	93130 UV	6.905	
22	18.19	0.00	1.00000	73653 VB	5.508	
23	22.42	0.00	1.00000	36212 BV	2.858	
24	23.47	0.00	1.00000	40730 UV	3.047	
25	27.12	0.00	1.00000	117044 VB	8.755	
26	35.55	0.00	1.00000	102307 BF	7.652	

TOTAL AREA = 1335917 TOTAL AREA % = 100.000

SAMPLE: P081016/1260 INJECTED AT 18:05:26 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 8 BTL: 8

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

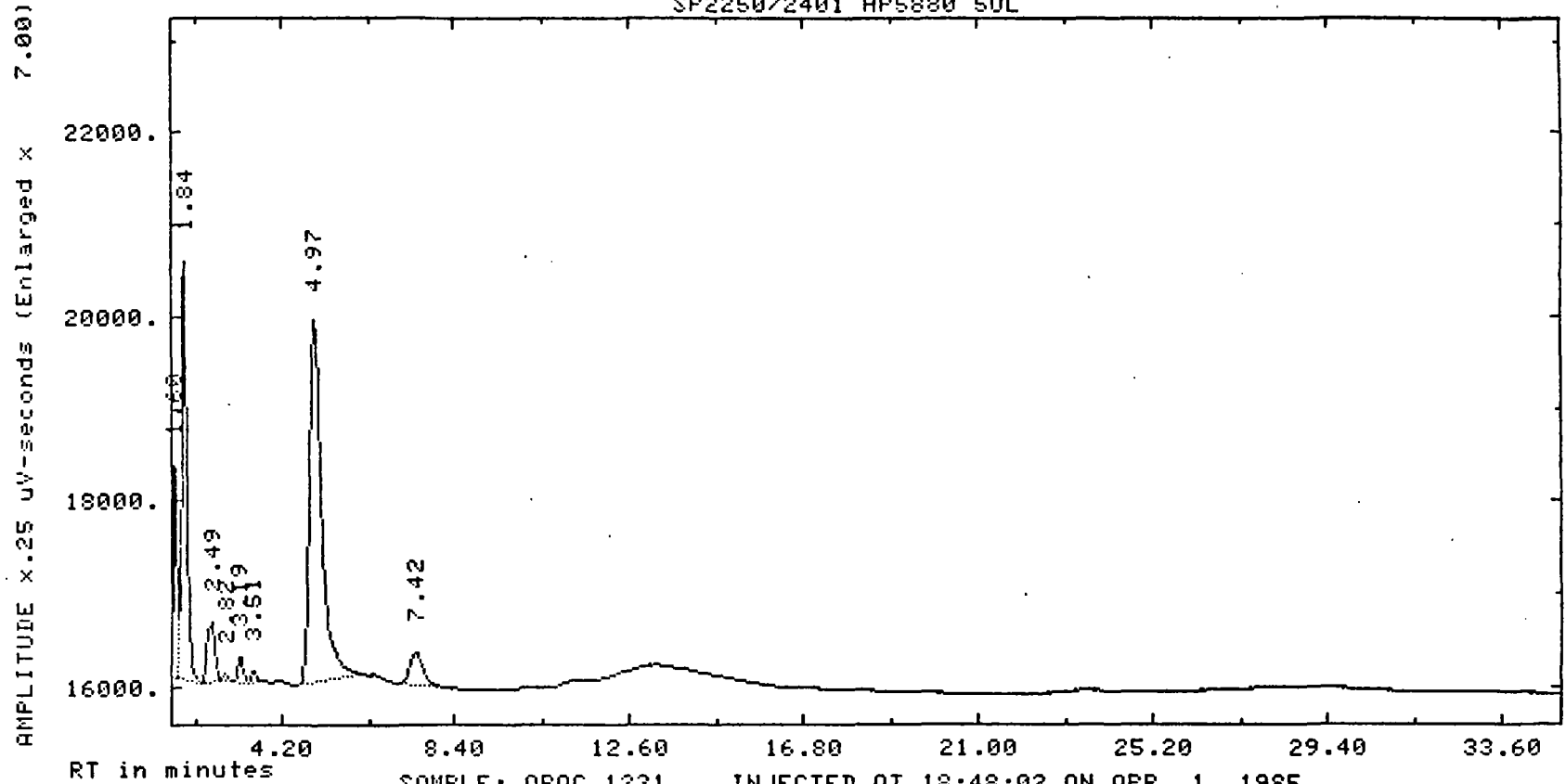
RAW DATA FILE: P85008 PARAM FILES= METHOD: SEQ:

DONE
READY

TV 158

Case 3969

SP2250/2401 HP5880 5UL



AMPLITUDE x.25 uV-seconds (Enlarged x 7.00)

RT in minutes

SAMPLE: AROC 1221 INJECTED AT 18:48:02 ON APR 1, 1985
Method: PRIME Raw: RB5009 Proc: PB5009

651 A

*LI,P,PS009

PROCESSED DATA FILE: PS009 ON CRN 21

Case 3969

APR 2, 1985 9:01:00

REPORT: 152 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HP5600 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.27	0.00	1.00000	1401 BV	1.756	
2	1.46	0.00	1.00000	1537 VV	1.924	
3	1.60	0.00	1.00000	7142 VV	8.945	
4	1.84	0.00	1.00000	19113 VB	23.938	
5	2.49	0.00	1.00000	4423 BV	5.539	
6	2.82	0.00	1.00000	300 VB	.376	
7	3.19	0.00	1.00000	1156 BV	1.461	
8	3.51	0.00	1.00000	497 VB	.623	
9	4.97	0.00	1.00000	40463 BB	50.701	
10	7.42	0.00	1.00000	3784 BB	4.740	

TOTAL AREA = 79846 TOTAL AREA % = 100.000

SAMPLE: AROC 1221 INJECTED AT 10:40:02 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 9 BTL: 9

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	.50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

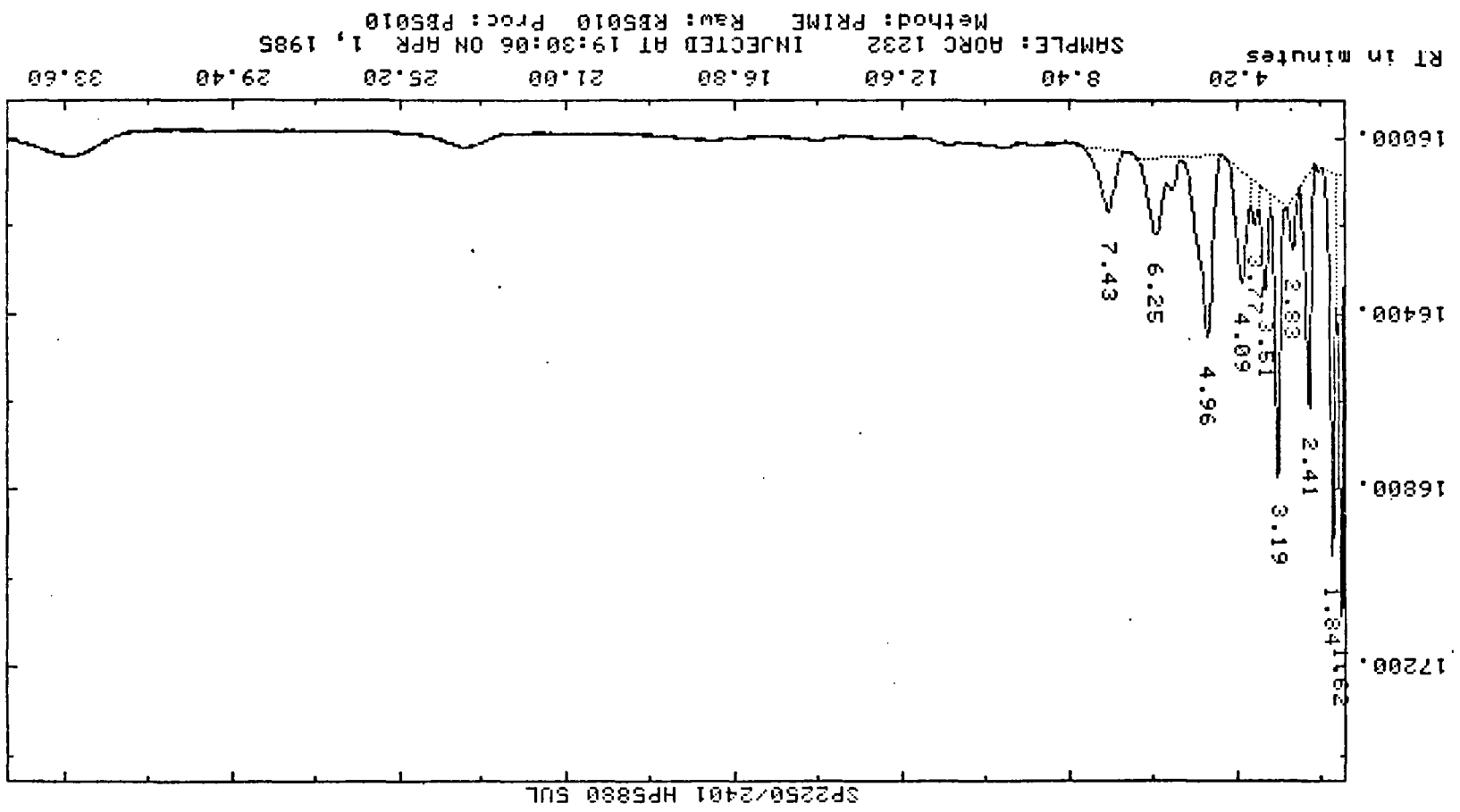
RE-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

RAW DATA FILE: PS009 PARAM FILES= METHOD: SEQ:

DONE

IV 160

AMPLITUDE X.25 UV-seconds (Enlarged X 150.00)



SP2250/2401 HP5880 SUL

Case 3969

IV 161

READY
*LI,P,P85010

PROCESSED DATA FILE: P85010 ON CRN 21

Case 3969

APR 2, 1985 9:11:05

REPORT: 153 CHANNEL: 2 # PEAKS: -13 SP2250/2401 HPS880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.27	0.00	1.00000	135 BV	.498	
2	1.46	0.00	1.00000	281 VV	.738	
3	1.62	0.00	1.00000	3781 VV	13.892	
4	1.64	0.00	1.00000	3454 VB	12.091	
5	2.41	0.00	1.00000	2578 BV	9.472	
6	2.83	0.00	1.00000	518 VB	1.875	
7	3.19	0.00	1.00000	2893 BV	10.628	
8	3.51	0.00	1.00000	1181 VV	4.340	
9	3.77	0.00	1.00000	588 VV	1.938	
10	4.09	0.00	1.00000	2281 VB	8.381	
11	4.95	0.00	1.00000	5095 BV	18.720	
12	5.25	0.00	1.00000	2687 VB	9.800	
13	7.43	0.00	1.00000	1940 BB	7.126	

TOTAL AREA = 27218 TOTAL AREA % = 100.000

SAMPLE: AORC 1232 INJECTED AT 19:30:08 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 10 BTL: 10

SL-WIDTH KV/MIN DELAY MIN-AR BUNCH REPORT
.500 .300 1.00 50 AUTO MEDIUM

SUP-UNK DVT ID-LVL REF-RTW %RTW %DIL-F ISO
NO 0.00 0 .30 5.0 100.00 YES

ACTUAL RUN TIME: 40.017 MINUTES

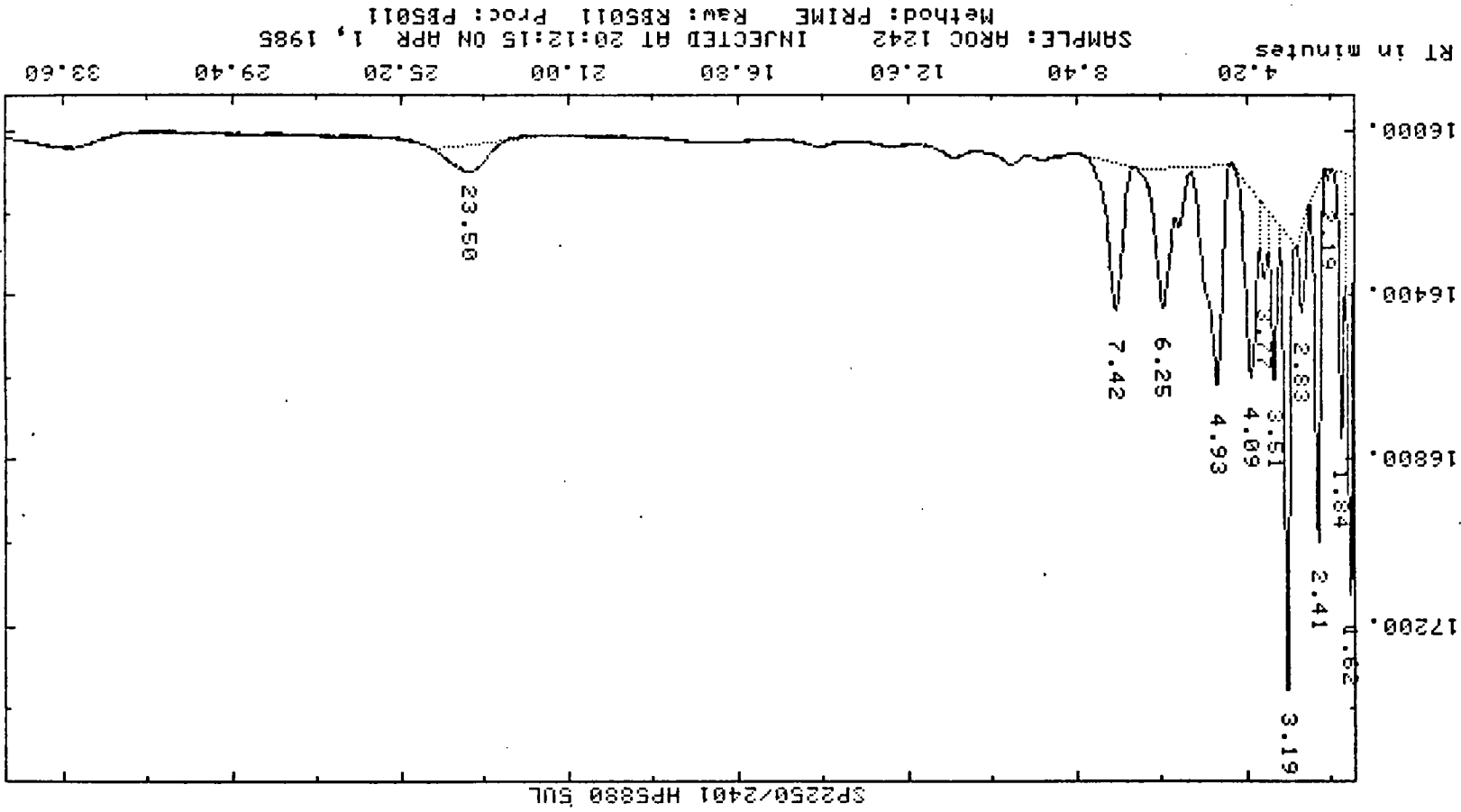
RF-UNK: 1.00000 STD-AMT: 1.00000 SAMP-AMT: 1.00000

RAW DATA FILE: P85010 PARAM FILES= METHOD: SEQ:

DONE
READY

TU 162

AMPLITUDE x.25 uV-seconds (Enlarged x 75.00)



Case 3969

IV 163

*L1,P,PB5011

PROCESSED DATA FILE: PB5011 ON CRN 21

Case 3969

APR 2, 1985 9:11:48

REPORT: 154 CHANNEL: 2 # PEAKS: -14 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.16	0.00	1.00000	87 BB	.204	
2	1.82	0.00	1.00000	4207 BV	9.918	
3	1.84	0.00	1.00000	2428 VB	5.668	
4	2.19	0.00	1.00000	61 BB	.142	
5	2.41	0.00	1.00000	3826 BV	8.931	
6	2.83	0.00	1.00000	945 VB	2.205	
7	3.19	0.00	1.00000	5271 BV	12.303	
8	3.51	0.00	1.00000	2862 VV	4.812	
9	3.77	0.00	1.00000	911 VV	2.126	
10	4.05	0.00	1.00000	4153 VV	9.094	
11	4.93	0.00	1.00000	7282 VV	16.996	
12	6.25	0.00	1.00000	5331 VB	12.444	
13	7.42	0.00	1.00000	4504 BB	10.513	
14	23.50	0.00	1.00000	1776 BB	4.146	

TOTAL AREA = 42844 TOTAL AREA % = 100.000

SAMPLE: AROC 1242 INJECTED AT 20:12:15 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: 5M05A SUBSQ/SAMP: 1 / 11 DTL: 11

SL-WDTH	NU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

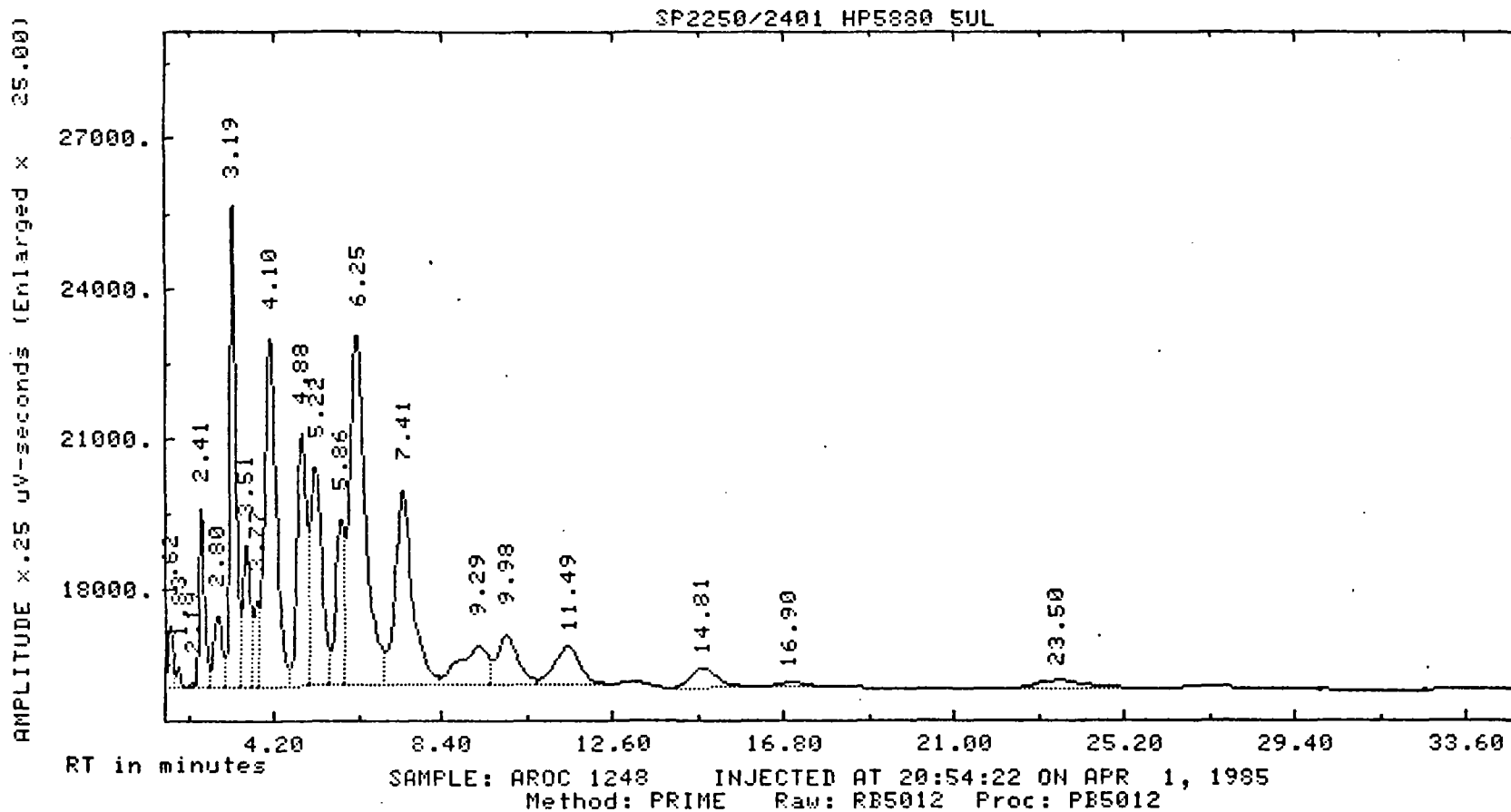
RAW DATA FILE: RBS011 PARAM FILES= METHOD: SEQ:

DONE
READY

10 164

Case 3969

591 AT



*LI,P,FB5012

PROCESSED DATA FILE: PD5012 ON CRN 21

Case 3969

APR 2, 1985 9:12:16

REPORT: 155 CHANNEL: 2 # PEAKS: -20 SP2250/2401 HP5880 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	5018 BV	.965	
2	1.83	0.00	1.00000	1350 VB	.260	
3	2.19	0.00	1.00000	276 BV	.653	
4	2.41	0.00	1.00000	16051 UV	3.088	
5	2.80	0.00	1.00000	9954 UV	1.921	
6	3.19	0.00	1.00000	51888 UV	9.950	
7	3.51	0.00	1.00000	17094 UV	3.440	
8	3.77	0.00	1.00000	7984 UV	1.535	
9	4.10	0.00	1.00000	70514 UV	13.557	
10	4.80	0.00	1.00000	41854 UV	8.050	
11	5.22	0.00	1.00000	41427 UV	7.954	
12	5.85	0.00	1.00000	26091 UV	5.016	
13	6.25	0.00	1.00000	95521 UV	18.575	
14	7.41	0.00	1.00000	62435 UV	12.003	
15	9.29	0.00	1.00000	18711 UV	3.597	
16	9.90	0.00	1.00000	17705 UV	3.404	
17	11.49	0.00	1.00000	16247 VB	3.124	
18	14.01	0.00	1.00000	9272 BV	1.783	
19	16.90	0.00	1.00000	2173 VB	.418	
20	23.50	0.00	1.00000	6601 BV	1.289	

TOTAL AREA = 520145 TOTAL AREA % = 100.000

SAMPLE: AROC 1240 INJECTED AT 20:54:22 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 12 BTL: 12

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.025 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

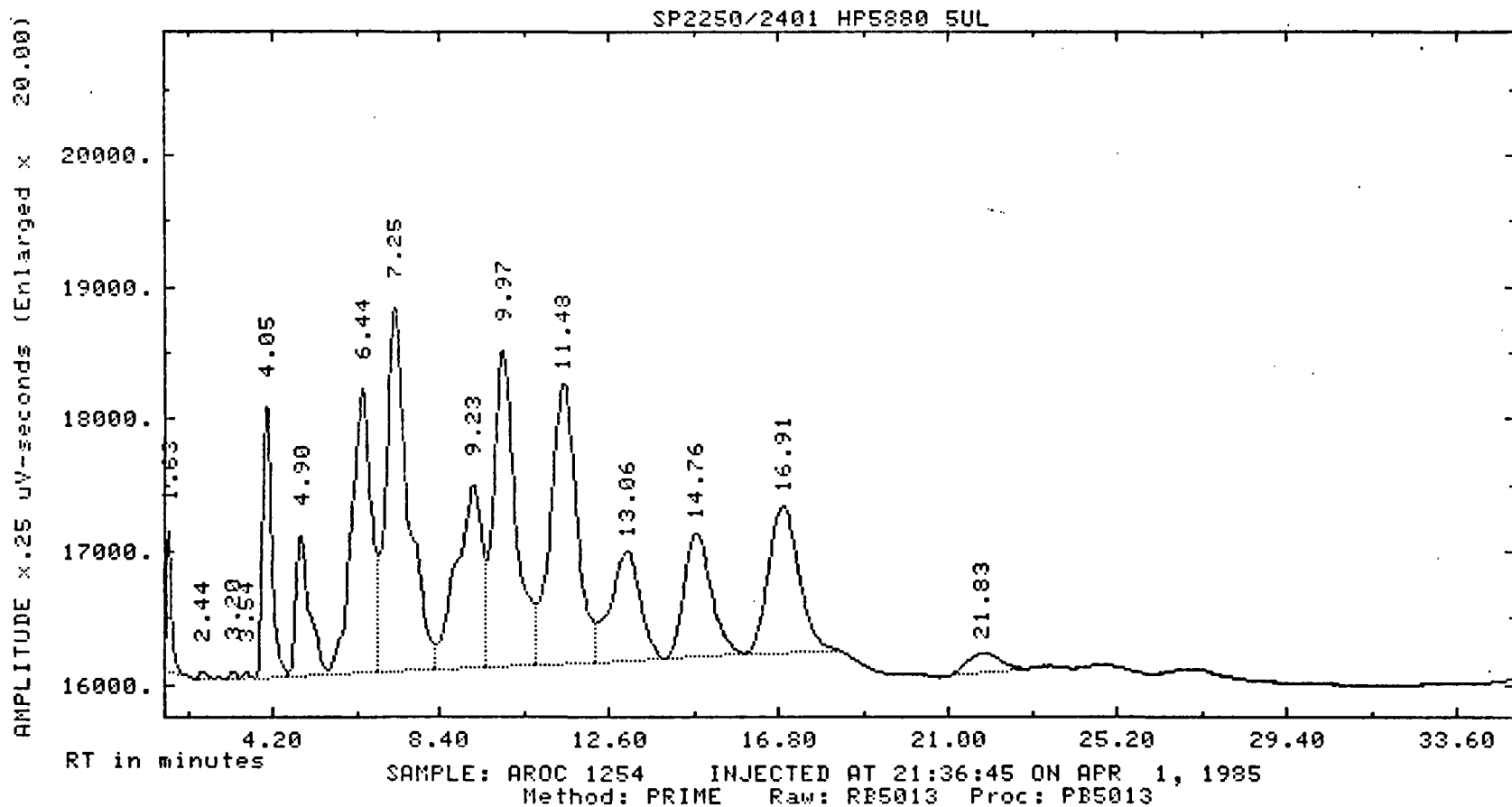
RAW DATA FILE: PD5012 PARAM FILES= METHOD: SEQ:

DONE
READY

Tv 166

Case 3969

L91 II



*LI,P,PB5013.

PROCESSED DATA FILE: PB5013 ON CRN 21

Case 3969

APR 2, 1985 9:12:47

REPORT: 156 CHANNEL: 2 # PEAKS: -16 SP2250/2401 HP5800 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.16	0.00	1.00000	109 EB	.033	
2	1.63	0.00	1.00000	4606 EB	1.411	
3	2.44	0.00	1.00000	281 EB	.086	
4	3.20	0.00	1.00000	275 EB	.084	
5	3.54	0.00	1.00000	213 EB	.065	
6	4.05	0.00	1.00000	16000 BV	4.905	
7	4.50	0.00	1.00000	12447 VV	3.814	
8	6.44	0.00	1.00000	36540 VV	11.226	
9	7.25	0.00	1.00000	56377 VV	17.273	
10	9.23	0.00	1.00000	30547 VV	9.359	
11	9.97	0.00	1.00000	46151 VV	14.143	
12	11.40	0.00	1.00000	47562 VV	14.603	
13	13.06	0.00	1.00000	21117 VV	6.476	
14	14.76	0.00	1.00000	21548 VB	6.602	
15	16.91	0.00	1.00000	28208 EB	8.642	
16	21.83	0.00	1.00000	4192 EB	1.284	

TOTAL AREA - 326391 TOTAL AREA % - 100.000

SAMPLE: AROC 1254 INJECTED AT 21:36:45 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SH03A SUBSQ/SAMP: 1 / 13 BTL: 13

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

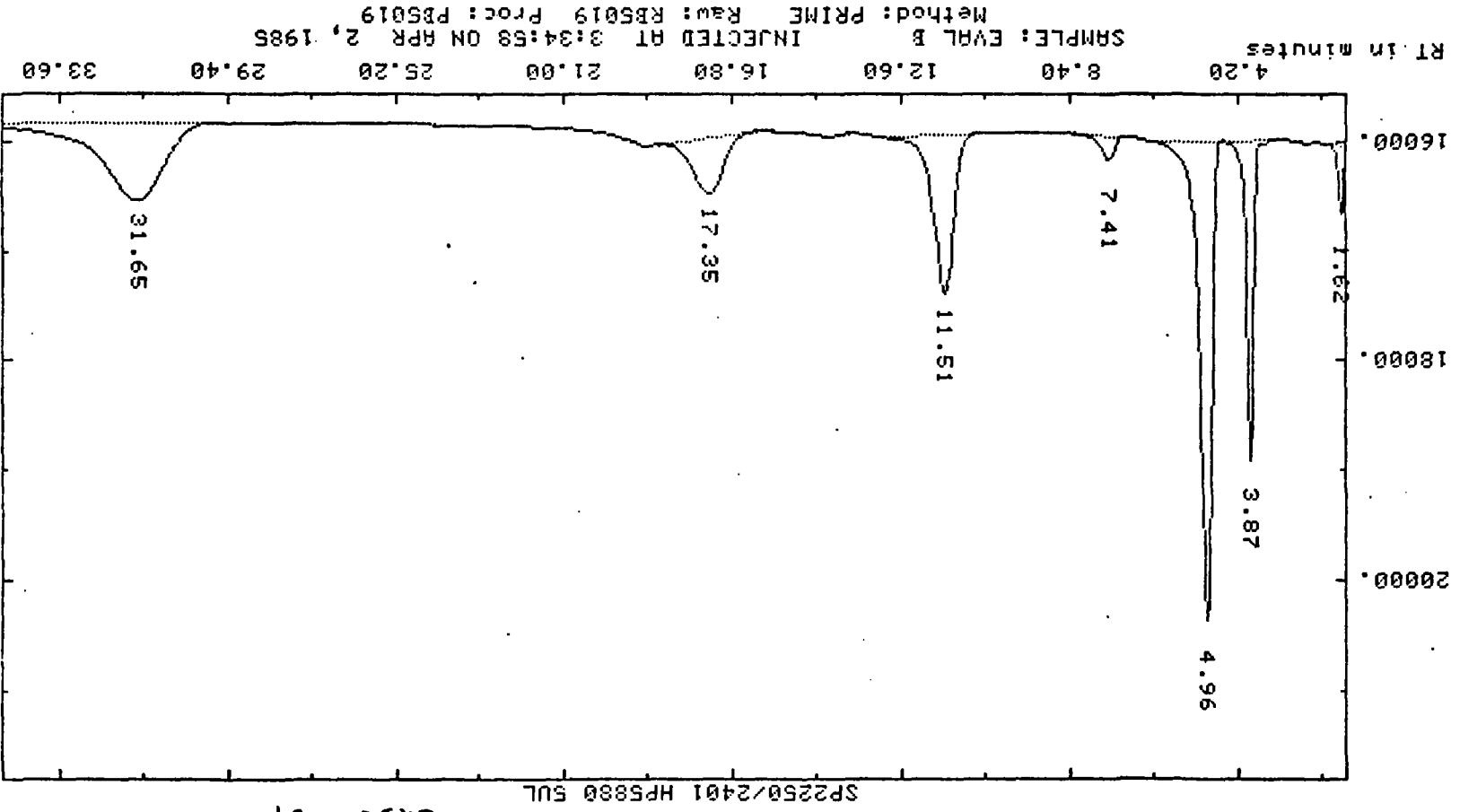
RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

RAW DATA FILE: PB5013 PARAM FILES= METHOD: SEQ:

DONE
READY

Jr 168

AMPLITUDE x.25 UV-seconds (Enlarged x 7.00)



SP2250/2401 HP5880 SUL

Case 3969

IV 169

*LI,P,PB5019

PROCESSED DATA FILE: PB5019 ON CRN 21

Case 3969

APR 2, 1985 9:16:09

REPORT: 162 CHANNEL: 2 # PEAKS: -8 SP2250/2401 HP5880 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.16	0.00	1.00000	112 BB	.076	
2	1.62	0.00	1.00000	2556 BB	1.719	
3	3.87	0.00	1.00000	18119 BB	12.185	Aldrin
4	4.96	0.00	1.00000	46246 BB	31.099	
5	7.41	0.00	1.00000	2585 BB	1.739	
6	11.51	0.00	1.00000	25195 BB	16.943	Endrin
7	17.35	0.00	1.00000	12764 BB	8.583	pp'-ODT
8	31.65	0.00	1.00000	41129 BB	27.650	Dibutylchlorodate

TOTAL AREA = 148707 TOTAL AREA % = 100.000

SAMPLE: EVAL B INJECTED AT 3:34:58 ON APR 2, 1985

ZERO METHOD: PRIKE SEQ: SMO3A SUBSQ/SAMP: 1 / 19 BTL: 19

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

RAW DATA FILE: PB5019 PARAM FILES= METHOD: SEQ:

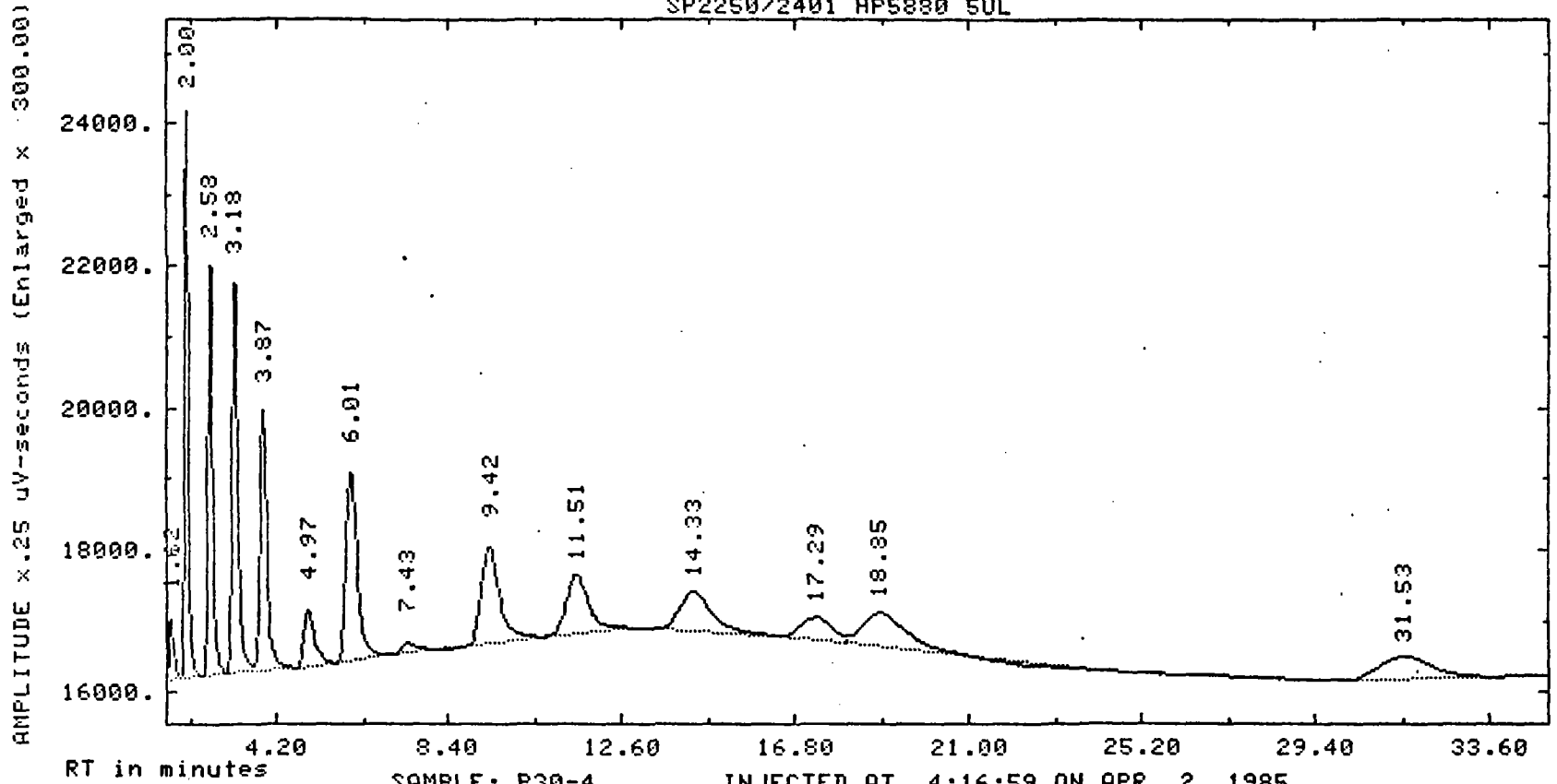
DONE
READY

IV 170

Case 3969

SP2250/2401 HP5880 5UL

141 II



SAMPLE: P30-4 INJECTED AT 4:16:59 ON APR 2, 1985
Method: PRIME Raw: RB5020 Proc: PB5020

*LI,P,PB5020

Case 3969

PROCESSED DATA FILE: PB5020 ON CRN 21

APR 2, 1985 9:16:33

REPORT: 163 CHANNEL: 2 # PEAKS: -14 SP2250/2401 HP5800 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	3417 BV	1.547	
2	2.00	0.00	1.00000	24762 VB	11.222	α -BHC
3	2.58	0.00	1.00000	22619 BV	10.243	γ -BHC
4	3.18	0.00	1.00000	28109 VV	12.729	Heptachlor
5	3.87	0.00	1.00000	21767 VB	9.857	Aldrin
6	4.97	0.00	1.00000	6920 BB	3.138	
7	6.01	0.00	1.00000	24305 BB	11.034	Heptachlor Epoxide
8	7.43	0.00	1.00000	1519 BB	.688	
9	9.42	0.00	1.00000	20266 BB	9.177	Dieldrin
10	11.51	0.00	1.00000	14705 BB	6.659	Endrin
11	14.33	0.00	1.00000	14541 BB	6.630	pp'-DDD
12	17.29	0.00	1.00000	8279 BV	3.749	pp'-DDT
13	18.05	0.00	1.00000	15721 VB	7.119	Endrin Aldehyde
14	31.53	0.00	1.00000	13706 BB	6.207	Pip Endrin Ketone

TOTAL AREA = 220824 TOTAL AREA % = 100.000

SAMPLE: P30-4 INJECTED AT 4:16:53 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SM03A SUBSQ/SAMP: 1 / 20 BTL: 20

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.000 MINUTES

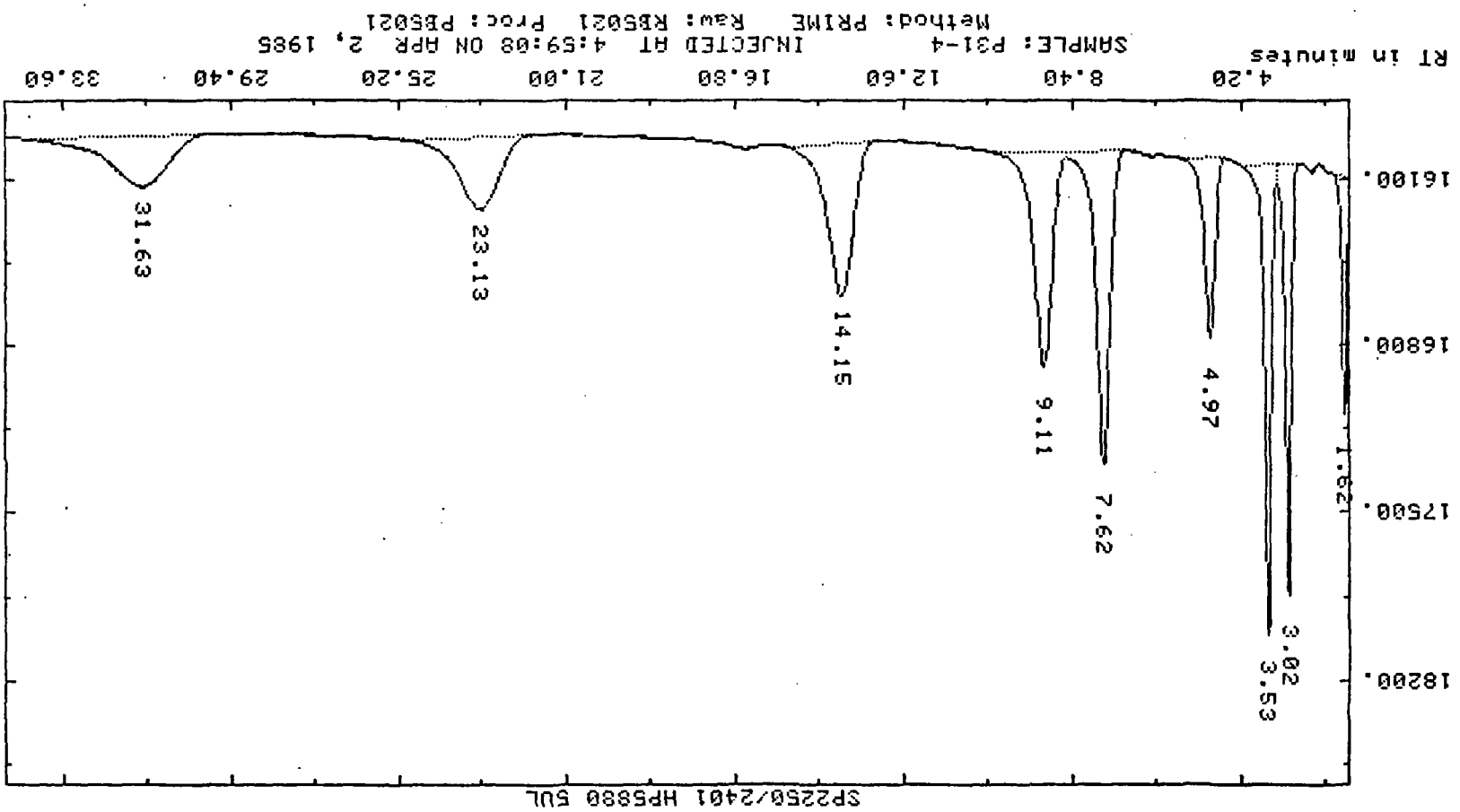
RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

RAW DATA FILE: RB5020 PARAM FILES= METHOD: SEQ:

DONE
READY

TV 172

AMPLITUDE x.25 UV-seconds (Enlarged x 25.00)



Case 3969

IV 173

SP2250/2401 HP5880 SUL

*LI,P,PB5021

PROCESSED DATA FILE: PB5021 ON CRN 21

Case 3969

APR 2, 1985 9:16:57

REPORT: 164 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HPS080 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.18	0.00	1.00000	119 BB	.119	
2	1.62	0.00	1.00000	4258 BB	4.280	
3	3.02	0.00	1.00000	9501 BV	9.526	B-BHC
4	3.53	0.00	1.00000	11692 VB	11.725	γ-BHC
5	4.97	0.00	1.00000	6957 EB	7.006	
6	7.62	0.00	1.00000	16210 BV	16.255	Endosulfan I
7	9.11	0.00	1.00000	14334 VB	14.374	PP'-DDE
8	14.15	0.00	1.00000	14292 EB	14.333	Endosulfan II
9	23.13	0.00	1.00000	11005 EB	11.035	
10	31.63	0.00	1.00000	11312 EB	11.344	Dibutyl chlorocate

TOTAL AREA = 99719 TOTAL AREA % = 100.000

SAMPLE: P31-4 INJECTED AT 4:59:08 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 21 BTL: 21

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.025 MINUTES

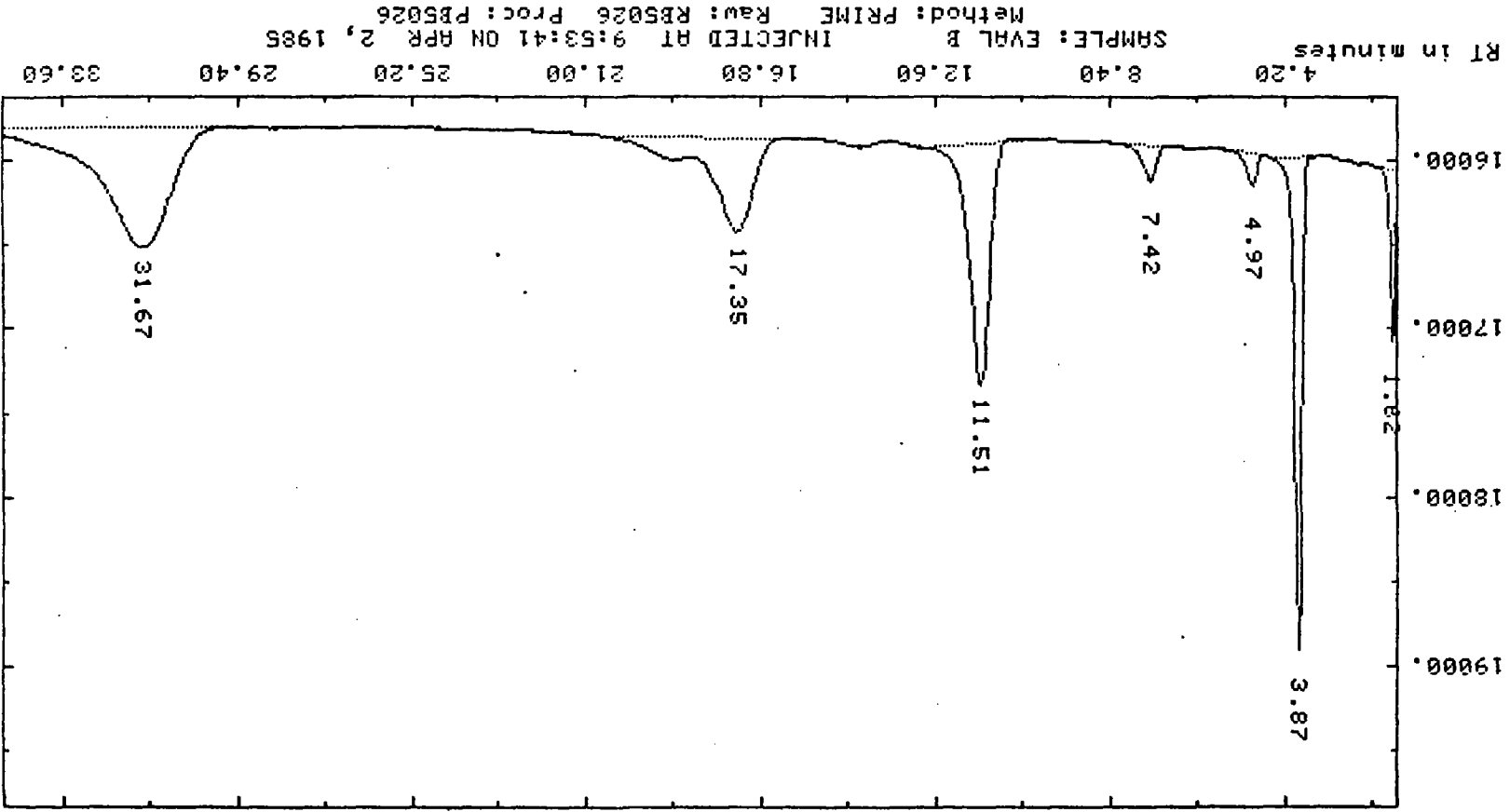
RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

RAW DATA FILE: PB5021 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 174

AMPLITUDE x.25 uV-seconds (Enlarged x 10.00)



SP2250/2401 HP5880 5UL

Case 3969

IV 175

Case 3969

LI, P, PB5026

PROCESSED DATA FILE: PB5020 ON CRN 21

APR 2, 1985 10:35:15

REPORT: 169 CHANNEL: 2 # PEAKS: -7 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	4282 BB	3.759	
2	3.87	0.00	1.00000	17653 BB	15.500	Aldrin
3	4.97	0.00	1.00000	1662 BB	1.460	
4	7.42	0.00	1.00000	2402 BB	2.105	
5	11.51	0.00	1.00000	24548 BB	21.554	Endrin
6	17.35	0.00	1.00000	20731 BB	18.203	pp'-DDT
7	31.67	0.00	1.00000	42612 BB	37.415	Dibutyltin diacetate

TOTAL AREA = 113891 TOTAL AREA % = 100.000

SAMPLE: EVAL B INJECTED AT 9:53:41 ON APR 2, 1985

ERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 26 BTL: 26

PL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

PUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISD
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

REP-UNK: 1.00000 STD-ANT: 1.0000 SAMP-ANT: 1.0000

RAW DATA FILE: PB5026 PARAM FILES= METHOD: SEQ:

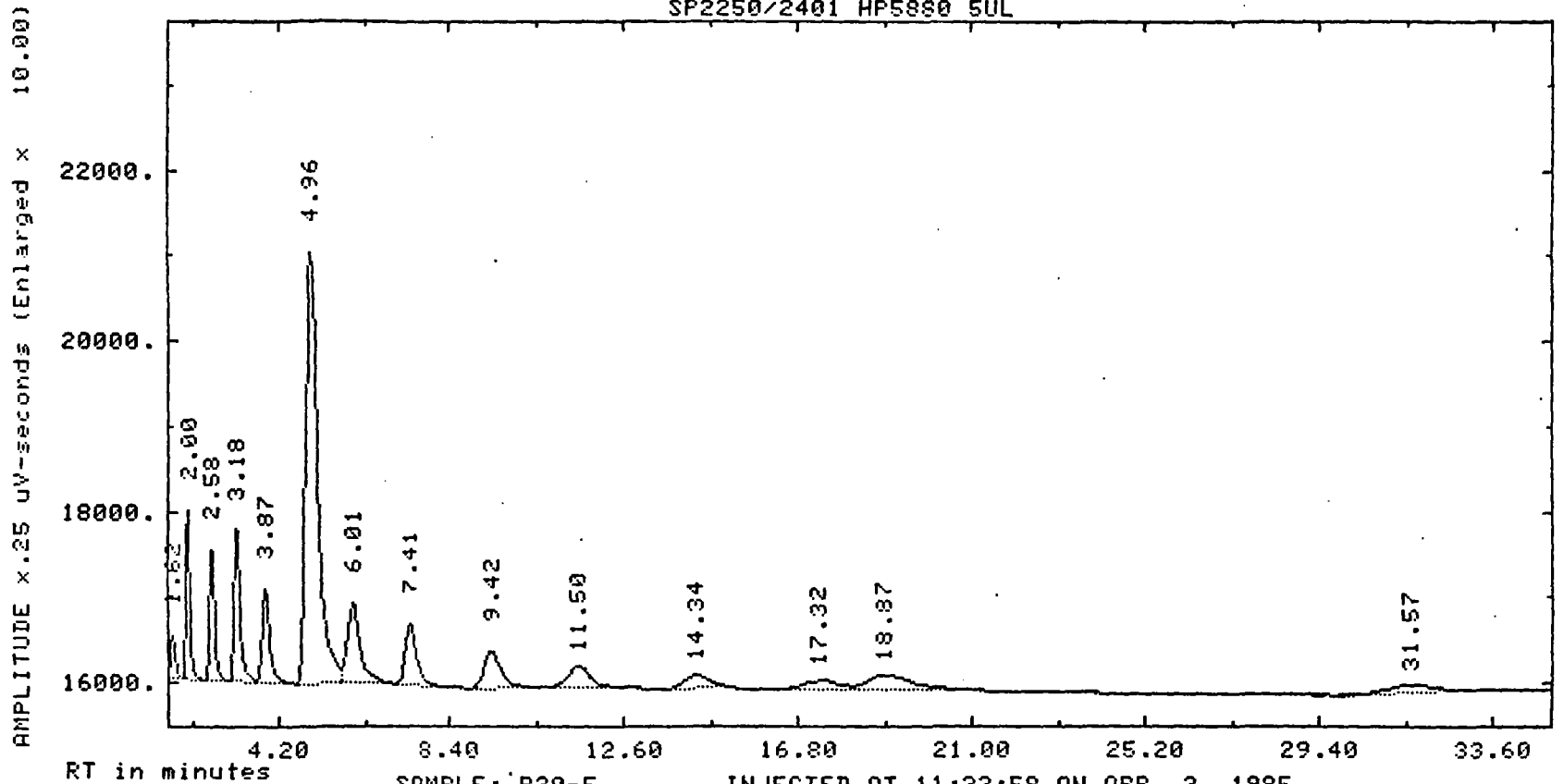
NONE

IV 176

Case 3969

SP2250/2401 HP5880 5UL

661 II



SAMPLE: P30-5 INJECTED AT 11:22:58 ON APR 2, 1985
Method: PRIME Raw: RB5027 Proc: PB5027

*LI,P,PD5027

PROCESSED DATA FILE: PB5027 ON CRN 21

Case 3969

APR 2, 1985 11:50:34

REPORT: 170 CHANNEL: 2 # PEAKS: -13 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.16	0.00	1.00000	.149 BB	.121	
2	1.63	0.00	1.00000	1884 BB	1.521	
3	2.00	0.00	1.00000	6955 BB	5.617	α -BHC
4	2.58	0.00	1.00000	6811 BV	5.501	γ -BHC
5	3.18	0.00	1.00000	9413 VB	7.602	Heptachlor
6	3.87	0.00	1.00000	6874 BB	5.551	Aldrin
7	4.97	0.00	1.00000	45860 BV	37.034	
8	6.01	0.00	1.00000	10041 VB	8.109	Heptachlor Epoxide
9	7.44	0.00	1.00000	2418 BB	1.953	
10	9.43	0.00	1.00000	6282 BB	6.688	Dieldrin
11	11.51	0.00	1.00000	3021 BB	6.478	Endrin
12	14.31	0.00	1.00000	12637 BB	10.205	pp'-DDO
13	31.46	0.00	1.00000	4483 BB	3.621	Endrin Ketone

TOTAL AREA = 123830 TOTAL AREA % = 100.000

SAMPLE: P30-5 INJECTED AT 10:35:39 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 27 BTL: 27

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.025 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

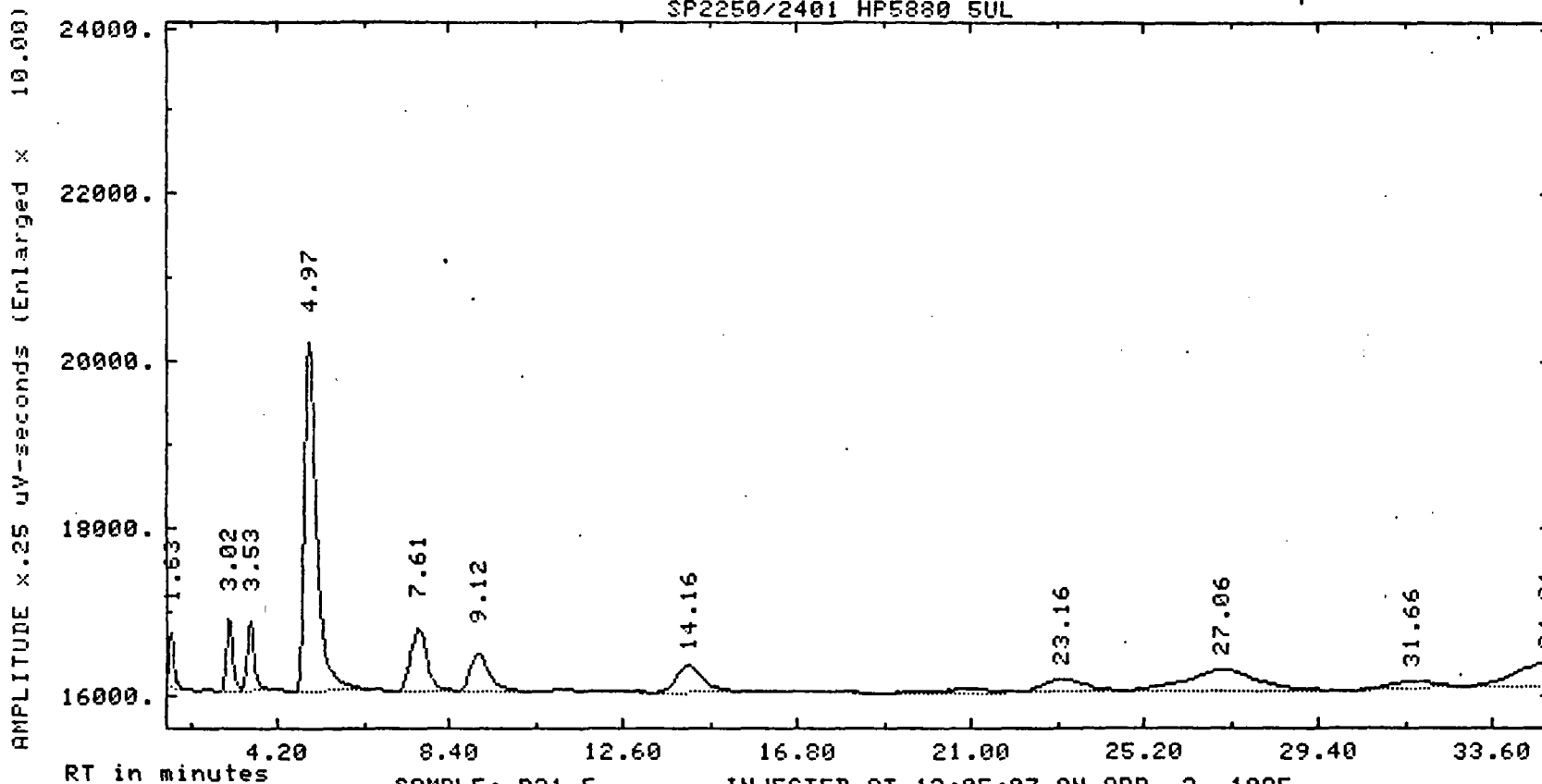
RAW DATA FILE: RBS027 PARAM FILES= METHOD: SEQ:

DONE

TU 178

Case 3969

SP2250/2401 HP5880 SUL



AMPLITUDE x.25 uV-seconds (Enlarged x 10.00)

RT in minutes

SAMPLE: P31-5 INJECTED AT 12:05:07 ON APR 2, 1985
Method: PRIME Raw: RB5028 Proc: PB5028

561 $\overline{17}$

93

READY
*LI,P,FB5028

Case 3969

PROCESSED DATA FILE: PB5028 ON CRN 21

APR 2, 1985 14:11:53

REPORT: 172 CHANNEL: 2 # PEAKS: -11 SP2250/2401 HP5890 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	2664 BB	2.240	
2	3.02	0.00	1.00000	4370 BV	3.675	B-BHC
3	3.53	0.00	1.00000	4817 VB	4.050	S-BHC
4	4.97	0.00	1.00000	42929 BB	36.098	
5	7.61	0.00	1.00000	10036 BB	8.438	Endosulfan I
6	9.12	0.00	1.00000	6545 BB	5.504	PP'-DDE
7	14.16	0.00	1.00000	6710 BB	5.642	Endosulfan II
8	23.16	0.00	1.00000	6035 BB	5.075	Endosulfan Sulfate
9	27.06	0.00	1.00000	13966 BB	11.744	
10	31.66	0.00	1.00000	2852 BB	2.398	Dibutylchloroendate
11	34.91	0.00	1.00000	18000 BB	15.136	

TOTAL AREA = 118925 TOTAL AREA % = 100.000

SAMPLE: P31-5 INJECTED AT 12:05:07 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 28 BTL: 28

SL-WIDTH MV/MIN DELAY MIN-AR BUNCH REPORT
.500 .300 1.00 50 AUTO MEDIUM

SUP-UNK DVT ID-LVL REF-RTW %RTW %DIL-F ISO
NO 0.00 0 .30 5.0 100.00 YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

RAW DATA FILE: RB5028 PARAM FILES= METHOD: SEQ:

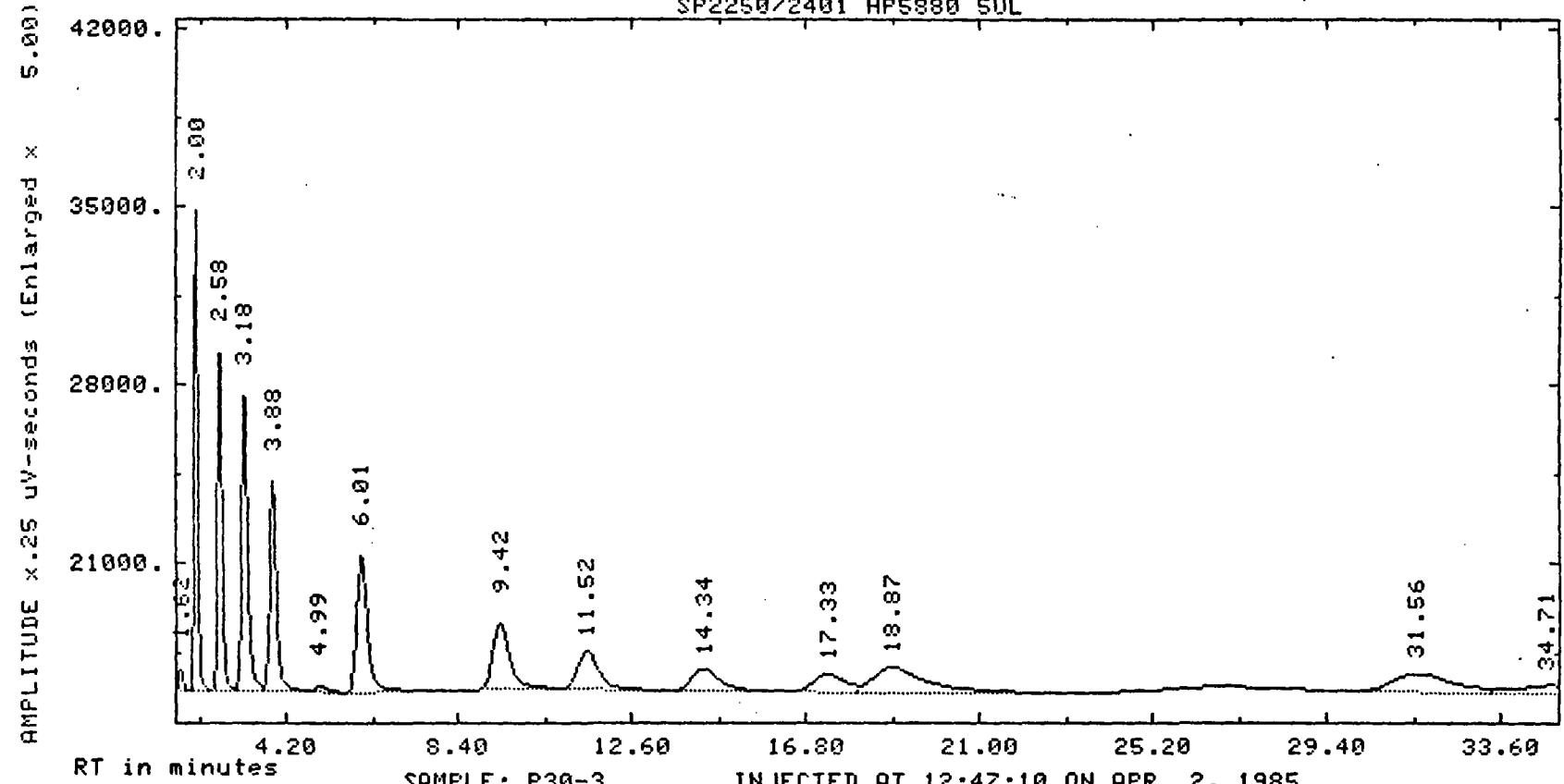
DONE
READY

IV 180

Case 3969

SP2250/2401 HP5980 5UL

181 II



SAMPLE: P30-3 INJECTED AT 12:47:10 ON APR 2, 1985
Method: PRIME Raw: RB5029 Proc: PB5029

*LI,P,PB5029

PROCESSED DATA FILE: PB5029 ON CRN 21

Case 3961

APR 2, 1985 14:12:19

REPORT: 173 CHANNEL: 2 # PEAKS: -15 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.16	0.00	1.00000	89 BB	.019	
2	1.62	0.00	1.00000	3404 BV	.728	
3	2.00	0.00	1.00000	58313 UV	12.473	α -BHC
4	2.58	0.00	1.00000	52120 UV	11.149	γ -BHC
5	3.18	0.00	1.00000	58295 UV	12.470	Heptachlor
6	3.80	0.00	1.00000	46493 VB	9.945	Alarin
7	4.99	0.00	1.00000	1685 BB	.350	
8	6.01	0.00	1.00000	48926 BB	10.465	Heptachlor Epoxide
9	9.42	0.00	1.00000	37070 BB	7.929	Dieldrin
10	11.52	0.00	1.00000	25483 BB	5.451	Endrin
11	14.34	0.00	1.00000	20910 BB	4.473	pp'-DDO
12	17.33	0.00	1.00000	18651 BV	3.990	pp'-DDT
13	18.87	0.00	1.00000	42418 VB	9.073	Endrin Aldehyde
14	31.56	0.00	1.00000	34765 BV	7.436	Endrin Ketone
15	34.71	0.00	1.00000	18875 VB	4.037	Methoxychlor

TOTAL AREA = 467456 TOTAL AREA % = 100.000

SAMPLE: P30-3 INJECTED AT 12:47:10 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 29 BTL: 29

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.025 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

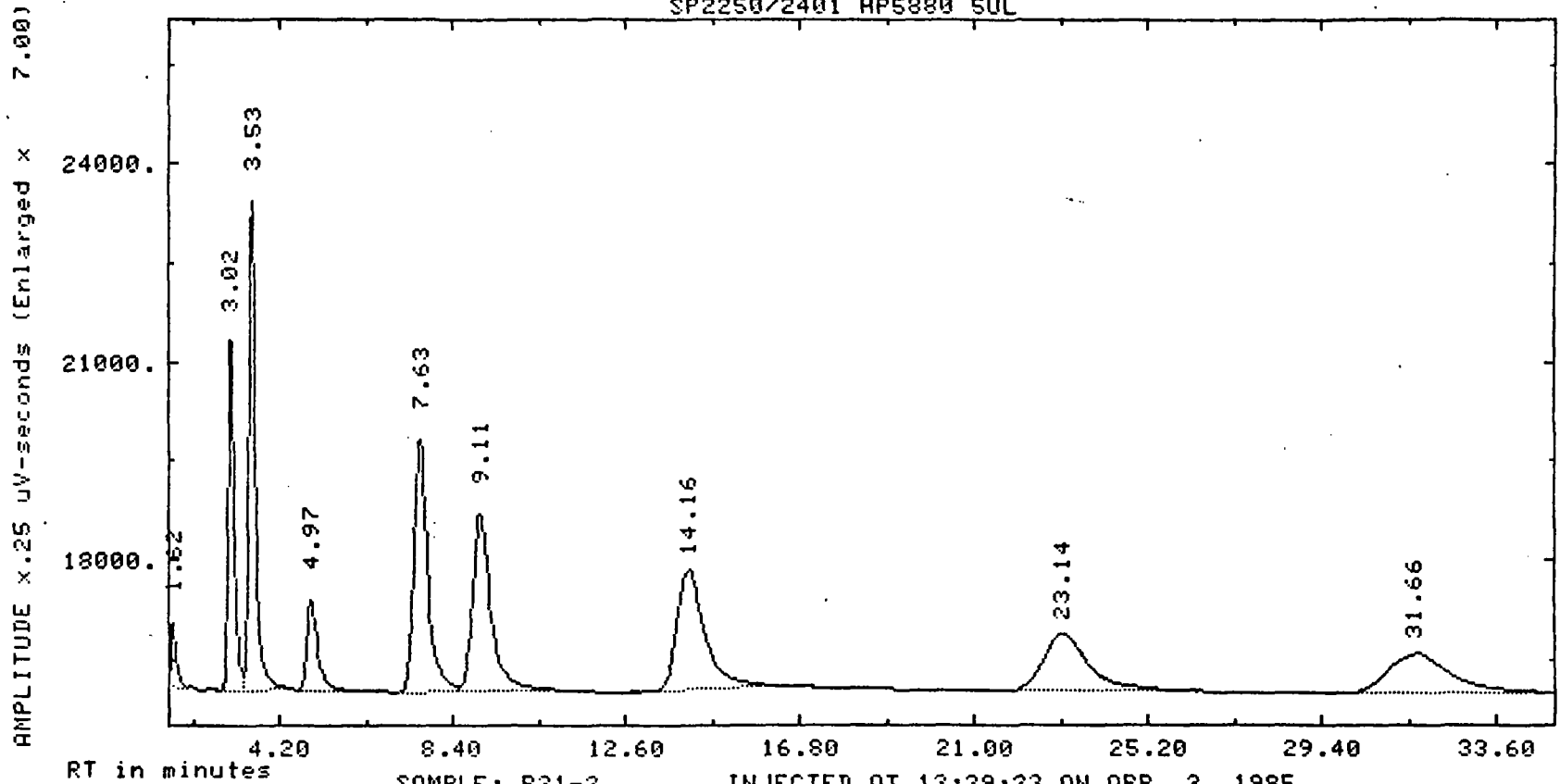
RAW DATA FILE: RB5029 PARAM FILES= METHOD: SEQ:

DONE
READY

10 182

Case 3969

SP2250/2401 HP5880 SUL



RT in minutes

SAMPLE: P31-3

INJECTED AT 13:29:23 ON APR 2, 1985

Method: PRIME Raw: RB5030 Proc: PB5030

IV 183

60

*LI,P,PB5030

PROCESSED DATA FILE: PB5030 ON CRN 21

Case 3969

APR 2, 1985 14:12:47

REPORT: 174 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HP5800 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.19	0.00	1.00000	135 BB	.049	
2	1.62	0.00	1.00000	4252 BB	1.538	
3	3.02	0.00	1.00000	27402 BV	9.910	B-BHC
4	3.53	0.00	1.00000	40629 VE	14.694	S-BHC
5	4.97	0.00	1.00000	12851 BB	4.648	
6	7.63	0.00	1.00000	45275 BV	16.374	Endosulfan I
7	9.11	0.00	1.00000	41393 VE	14.970	PP'-OPE
8	14.16	0.00	1.00000	39915 BB	14.436	Endosulfan II
9	23.14	0.00	1.00000	32244 BB	11.581	Endosulfan Sulfate
10	31.86	0.00	1.00000	32404 BB	11.719	Dibutylchloride

TOTAL AREA = 276501 TOTAL AREA % = 100.000

SAMPLE: P31-3 INJECTED AT 13:29:23 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 30 BTL: 30

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISQ
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.817 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-ANT: 1.0000

RAW DATA FILE: RE5030 PARAM FILES= METHOD: SEQ:

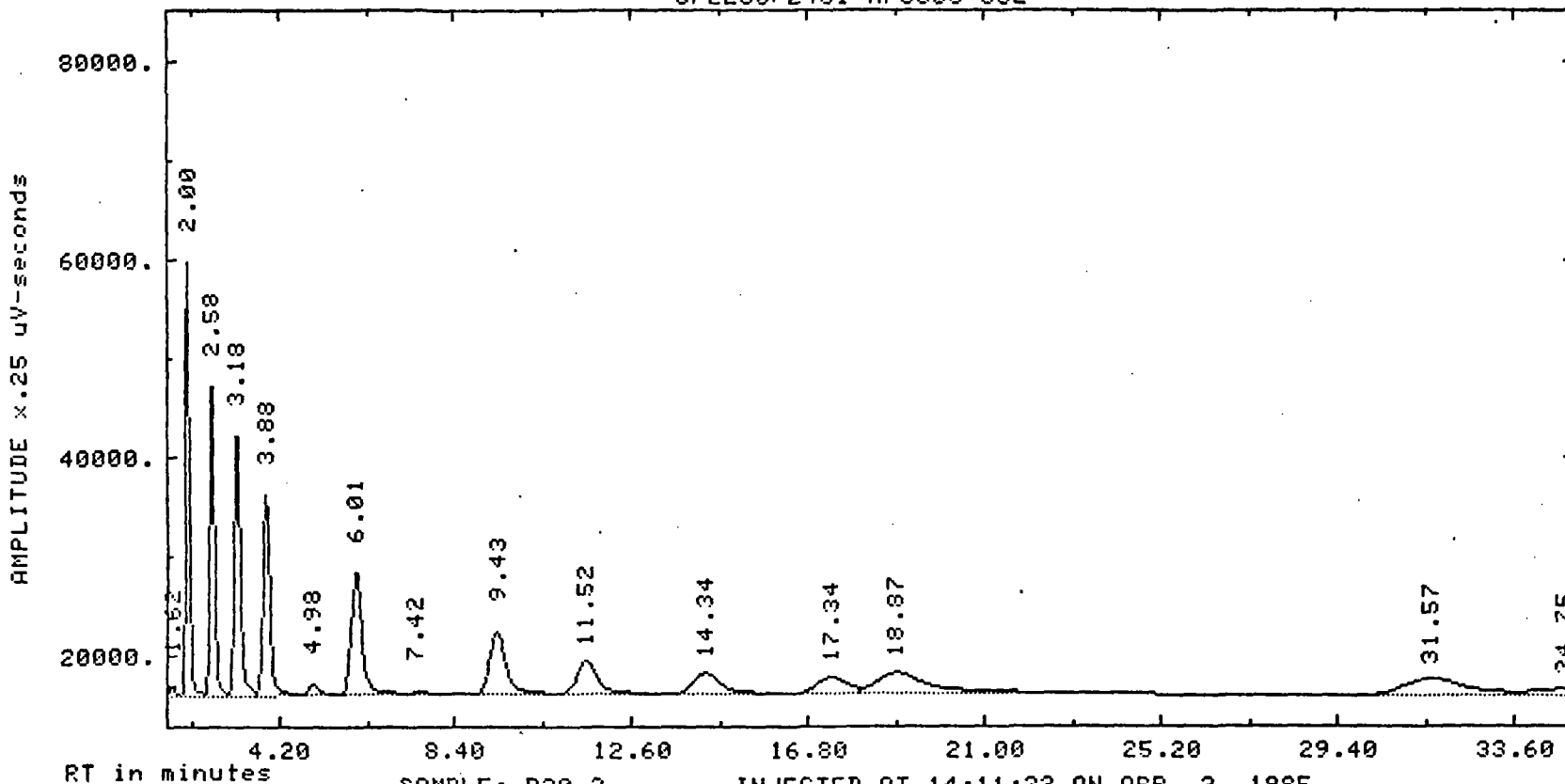
DONE

IV 184

Case 3969

SP2250/2401 HP5880 5UL

581 PT



SAMPLE: P30-2 INJECTED AT 14:11:23 ON APR 2, 1985
Method: PRIME Raw: RB5031 Proc: PB5031

READY
*LI,P,PE5031

PROCESSED DATA FILE: PE5031 ON CRN 21

Case 3969

APR 2, 1985 15:20:45

REPORT: 175 CHANNEL: 2 # PEAKS: -15 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.62	0.00	1.00000	3989 BV	.357	
2	2.08	0.00	1.00000	145409 UV	13.006	<i>a-BHC</i>
3	2.58	0.00	1.00000	130008 UV	11.628	<i>γ-BHC</i>
4	3.18	0.00	1.00000	134595 UV	12.039	<i>Heptachlor</i>
5	3.88	0.00	1.00000	117560 UV	10.515	<i>Aldrin</i>
6	4.95	0.00	1.00000	9743 UV	.871	
7	6.01	0.00	1.00000	113601 BV	10.161	<i>Heptachlor Epoxide</i>
8	7.42	0.00	1.00000	3014 UV	.270	
9	9.43	0.00	1.00000	87258 BV	7.805	<i>Dieldrin</i>
10	11.52	0.00	1.00000	59783 UV	5.347	<i>Endrin</i>
11	14.34	0.00	1.00000	50437 BV	4.511	<i>PP'-DDD</i>
12	17.34	0.00	1.00000	44722 UV	4.000	<i>PP'-DDT</i>
13	18.87	0.00	1.00000	94166 UV	8.422	<i>Endrin Alkylate</i>
14	31.57	0.00	1.00000	75900 BV	7.147	<i>Endrin Ketone</i>
15	34.75	0.00	1.00000	43240 UV	3.921	<i>Methoxychlor</i>

TOTAL AREA = 1118035 TOTAL AREA % = 100.000

SAMPLE: P30-2 INJECTED AT 14:11:23 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 31 BTL: 31

SL-WIDTH HV/MIN DELAY MIN-AR BUNCH REPORT
.500 .300 1.00 50 AUTO MEDIUM

SUP-UNK DVT ID-LVL REF-RTW %RTW %DIL-F ISO
NO 0.00 0 .30 5.0 100.00 YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

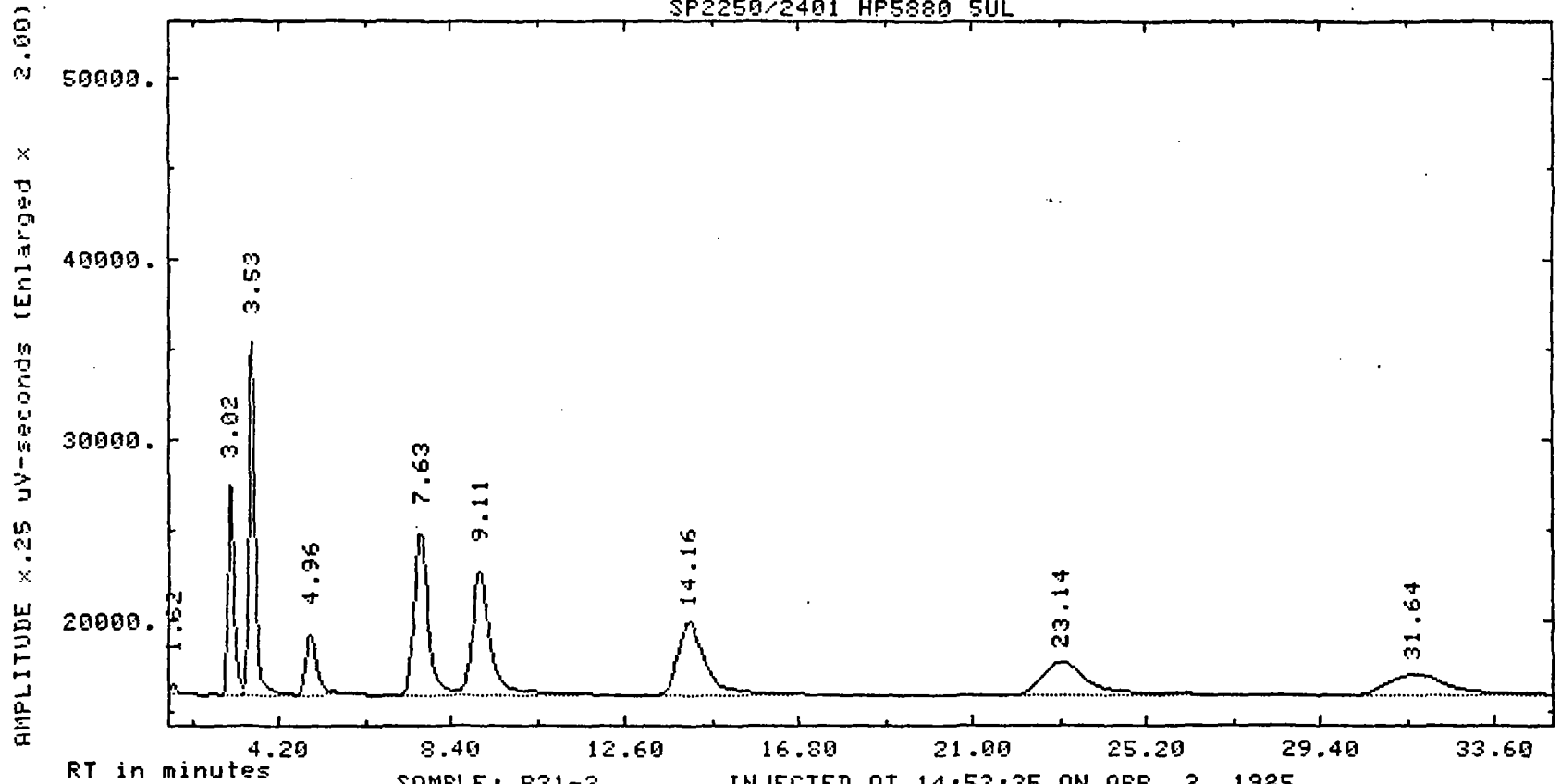
RAW DATA FILE: PE5031 PARAM FILES= METHOD: SEQ:

DONE

IV 186

Case 3969

SP2250/2401 HP5880 SUL



RT in minutes

SAMPLE: P31-2

INJECTED AT 14:53:35 ON APR 2, 1985

Method: PRIME Raw: RB5032 Proc: PB5032

LSI III

9.11

LI.P, PB5032

PROCESSED DATA FILE: PB5032 ON CRN 21

Case 3769

APR 2, 1985 15:35:01

REPORT: 176 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HP5980 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.19	0.00	1.00000	119 DB	.019	
2	1.62	0.00	1.00000	2388 BB	.379	
3	3.02	0.00	1.00000	50432 BV	9.596	B-BHC
4	3.53	0.00	1.00000	108013 UV	17.152	S-BHC
5	4.96	0.00	1.00000	32439 VB	5.151	
6	7.63	0.00	1.00000	104343 BV	16.570	Endosulfan I
7	9.11	0.00	1.00000	101215 VB	16.073	pp'-DDE
8	14.16	0.00	1.00000	91582 BD	14.543	Endosulfan II
9	23.14	0.00	1.00000	69502 EB	11.037	Endosulfan Sulfate
10	31.64	0.00	1.00000	59695 BB	9.479	Dib.tylchlorodete

TOTAL AREA = 629729 TOTAL AREA % = 100.000

SAMPLE: P31-2 INJECTED AT 14:53:35 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 32 BTL: 32

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	XRTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-ANT: 1.0000 SAMP-AMT: 1.0000

RAW DATA FILE: RES032 PARAM FILES= METHOD: SEQ:

DONE

IV 188

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2.0.
3.19

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5.01

2.45

21.14

2.207

2.231

2.91

Case 3969

SP2250/2401 HP5880 SUL

11-1
19

READY
READY
PLI.P.PB5033

PROCESSED DATA FILE: PD5033 ON CRN 21

Case 3969

APR 2, 1985 17:13:14

REPORT: 177 CHANNEL: 2 # PEAKS: -17 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.04	0.00	1.00000	59 BB	.003	
2	1.32	0.00	1.00000	483 BV	.026	
3	1.62	0.00	1.00000	2769 UV	.147	
4	2.01	0.00	1.00000	246057 UV	13.100	α -BHC
5	2.59	0.00	1.00000	220223 UV	11.724	δ -BHC
6	3.19	0.00	1.00000	227903 UV	12.133	Heptachlor
7	3.89	0.00	1.00000	217375 UV	11.573	Aldrin
8	5.01	0.00	1.00000	17843 UV	.950	
9	6.02	0.00	1.00000	187601 UV	9.987	Heptachlor Epoxide
10	7.43	0.00	1.00000	8307 VB	.442	
11	9.44	0.00	1.00000	162290 EV	8.640	Dieldrin
12	11.54	0.00	1.00000	118712 UV	6.320	Endrin
13	14.37	0.00	1.00000	77034 UV	4.101	PP'-DDD
14	17.30	0.00	1.00000	69893 UV	3.721	PP'-DDT
15	18.91	0.00	1.00000	140659 VB	7.489	Endrin Aldehyde
16	31.69	0.00	1.00000	121022 DV	6.443	Endrin Ketone
17	34.90	0.00	1.00000	60126 VB	3.201	Methoxychlor

TOTAL AREA = 1878357 TOTAL AREA % = 100.000

SAMPLE: P30-1 INJECTED AT 15:35:36 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 33 BTL: 33

SL-WIDTH MV/MIN DELAY -MIN-AR BUNCH REPORT
.500 .300 1.00 50 AUTO MEDIUM

SUP-UNK DVT ID-LVL REF-RTW %RTW %DIL-F ISO
NO 0.00 0 .30 5.0 100.00 YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

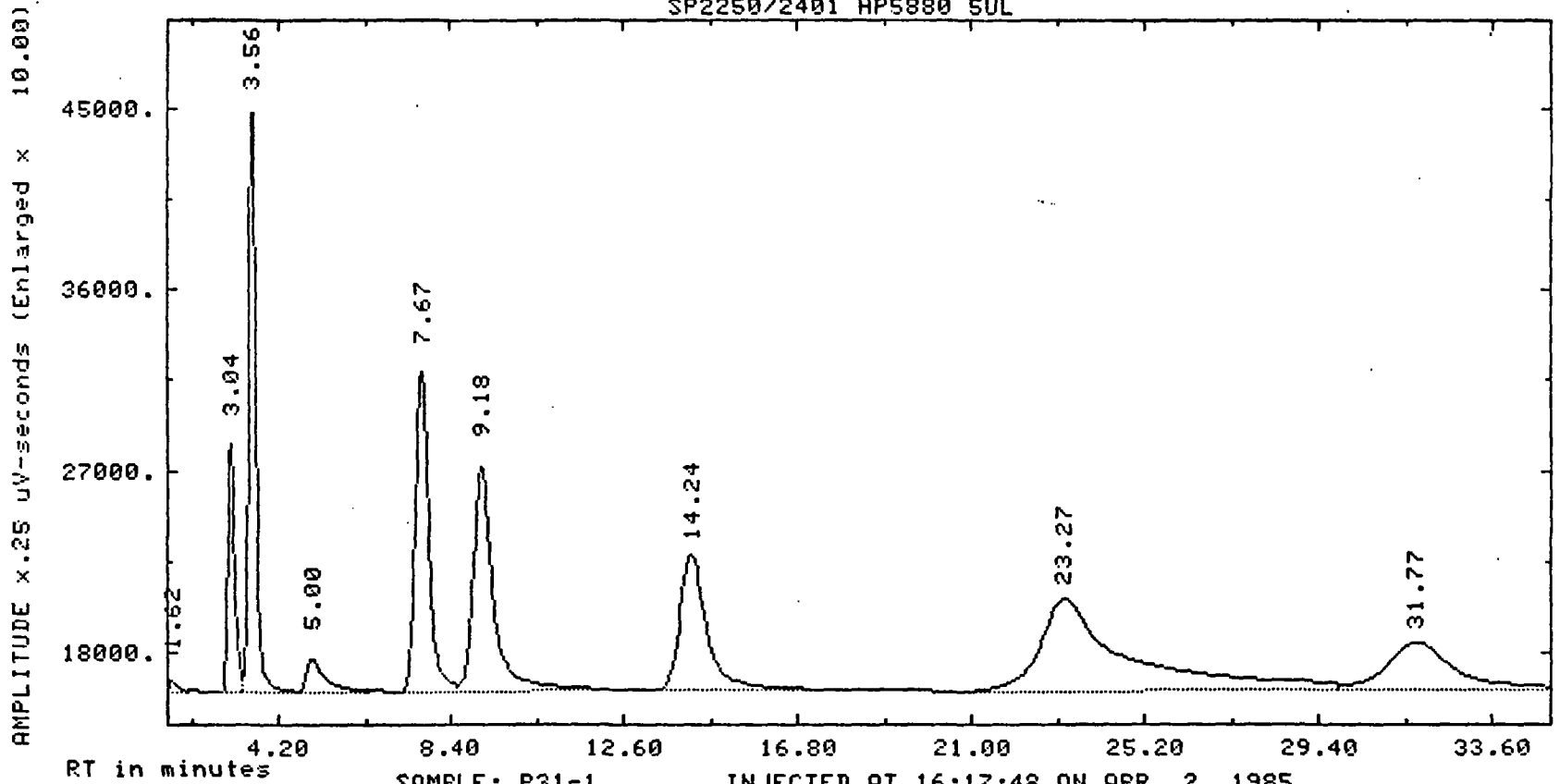
RAW DATA FILE: R95033 PARAM FILES= METHOD: SEQ:

DONE
READY

IV 190

Case 3969

SP2250/2401 HP5880 5UL



AMPLITUDE x.25 uV-seconds (Enlarged x 10.00)

RT in minutes

SAMPLE: P31-1

INJECTED AT 16:17:48 ON APR 2, 1985

Method: PRIME

Raw: RB5034

Proc: PB5034

161 PT

dv

*LI,P,PE5034

PROCESSED DATA FILE: PE5034 ON CRN 21

Case 3969

APR 2, 1995 17:13:39

REPORT: 178 CHANNEL: 2 # PEAKS: -10 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.20	0.00	1.00000	98 BB	.008	
2	1.62	0.00	1.00000	2920 BB	.231	
3	3.04	0.00	1.00000	72813 BV	5.764	B-BHC
4	3.56	0.00	1.00000	177697 VV	14.066	δ-BHC
5	5.00	0.00	1.00000	24182 VD	1.914	
6	7.67	0.00	1.00000	183237 BV	14.505	Endosulfan I
7	9.18	0.00	1.00000	189000 VB	14.969	PP'-DDE
8	14.24	0.00	1.00000	152702 BB	12.094	Endosulfan II
9	23.27	0.00	1.00000	327855 BV	25.954	Endosulfan Sulfate
10	31.77	0.00	1.00000	132582 VB	10.495	Dibutylchlorodate

TOTAL AREA = 1263274 TOTAL AREA % = 100.000

SAMPLE: P31-1 INJECTED AT 16:17:49 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMOZA SUBSQ/SAMP: 1 / 34 BTL: 34

SL-WDTH	MU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

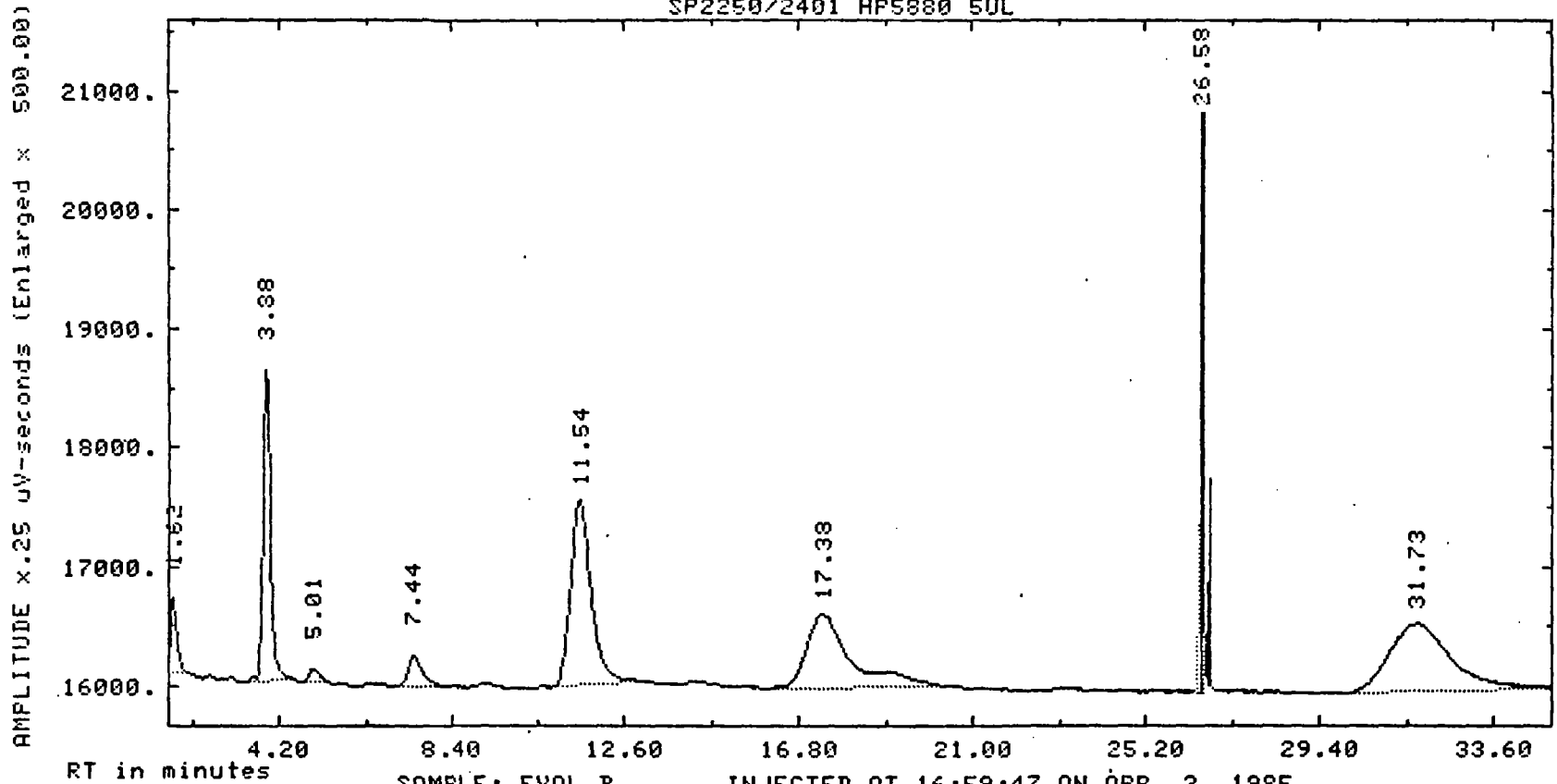
RAW DATA FILE: PE5034 PARAN FILES= METHOD: SEQ:

DONE

TV 192

Case 3969

SP2250/2401 HP5880 5UL



AMPLITUDE x.25 uV-seconds (Enlarged x 500.00)

RT in minutes

SAMPLE: EVAL B INJECTED AT 16:59:47 ON APR 2, 1985
Method: PRIME Raw: RB5035 Proc: PB5035

531 11

READY
-LI.P.P85035

PROCESSED DATA FILE: P85035 ON CRN 21

Case 3969

APR 3, 1985 10:02:26

REPORT: 179 CHANNEL: 2 # PEAKS: -9 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.31	0.00	1.00000	180 BB	.008	
2	1.62	0.00	1.00000	3174 BB	.141	
3	3.98	0.00	1.00000	16367 BB	.725	Aldrin
4	5.01	0.00	1.00000	902 BB	.040	
5	7.44	0.00	1.00000	3198 BB	.142	
6	11.54	0.00	1.00000	26674 BB	1.182	Endrin
7	17.38	0.00	1.00000	22181 BB	.982	PP'-DDT
8	26.58	0.00	1.00000	2154482 BS	95.431	
9	31.73	0.00	1.00000	30477 BB	1.350	Dibutylchlorohydrate

TOTAL AREA = 2257637 TOTAL AREA % = 100.000

SAMPLE: EVAL B INJECTED AT 16:59:47 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 35 BTL: 35

SL-WDTH MV/MIN DELAY MIN-AR BUNCH REPORT
.500 .300 1.00 50 AUTO MEDIUM

SUP-UNK DVT ID-LVL REF-RTW %RTW %DIL-F ISO
NO 0.00 0 .30 5.0 100.00 YES

ACTUAL RUN TIME: 40.017 MINUTES

RE-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

RAW DATA FILE: R85035 PARAM FILES= METHOD: SEQ:

DONE

TV 194

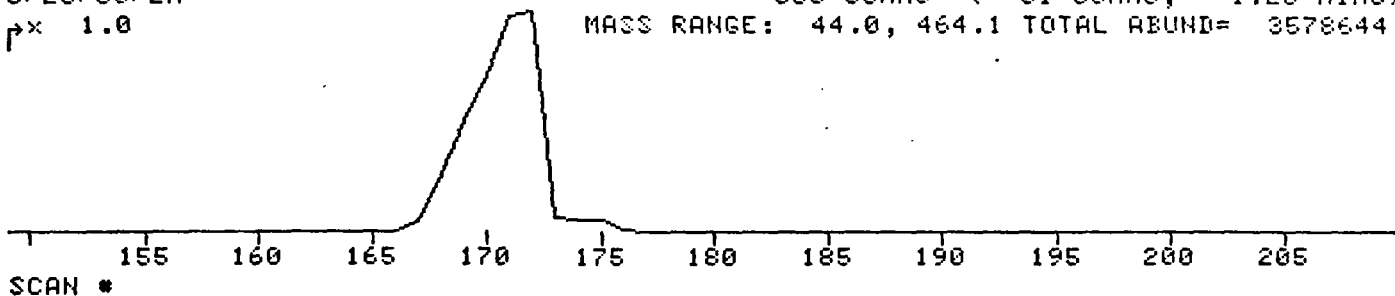
V. K&W QC DATA PACKET



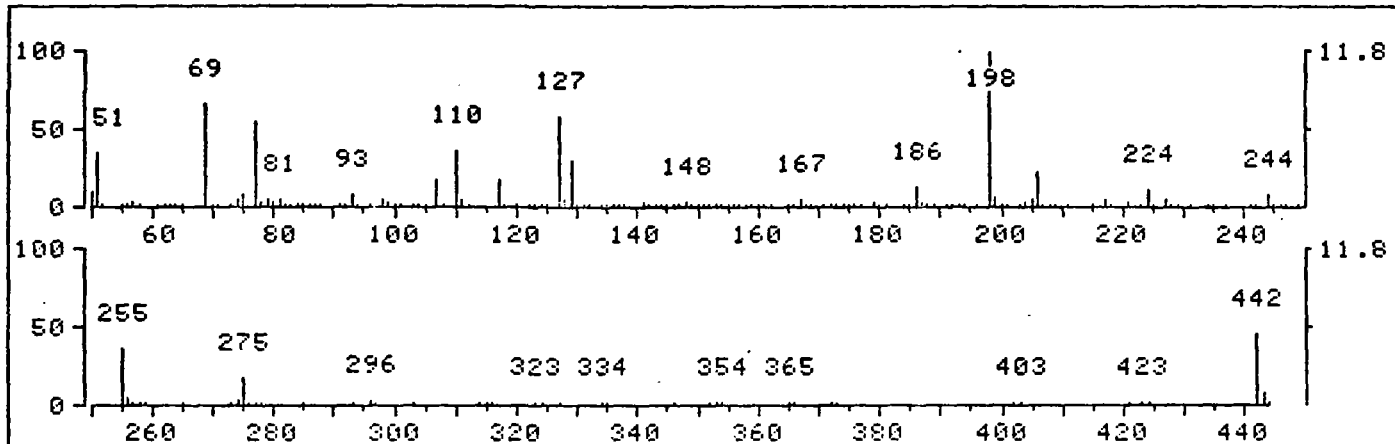
GCA CORPORATION
Technology Division

DFTPP
3/25/85/EM
p x 1.0

383 SCANS (61 SCANS, 1.28 MINS)
MASS RANGE: 44.0, 464.1 TOTAL ABUND= 3578644.



* 172 RET. TIME: 5.60 TOT ABUND= 222234. BASE PK/ABUND: 198.0/ 26314.



J-1

DRN: 14741
SPECTRUM: 172

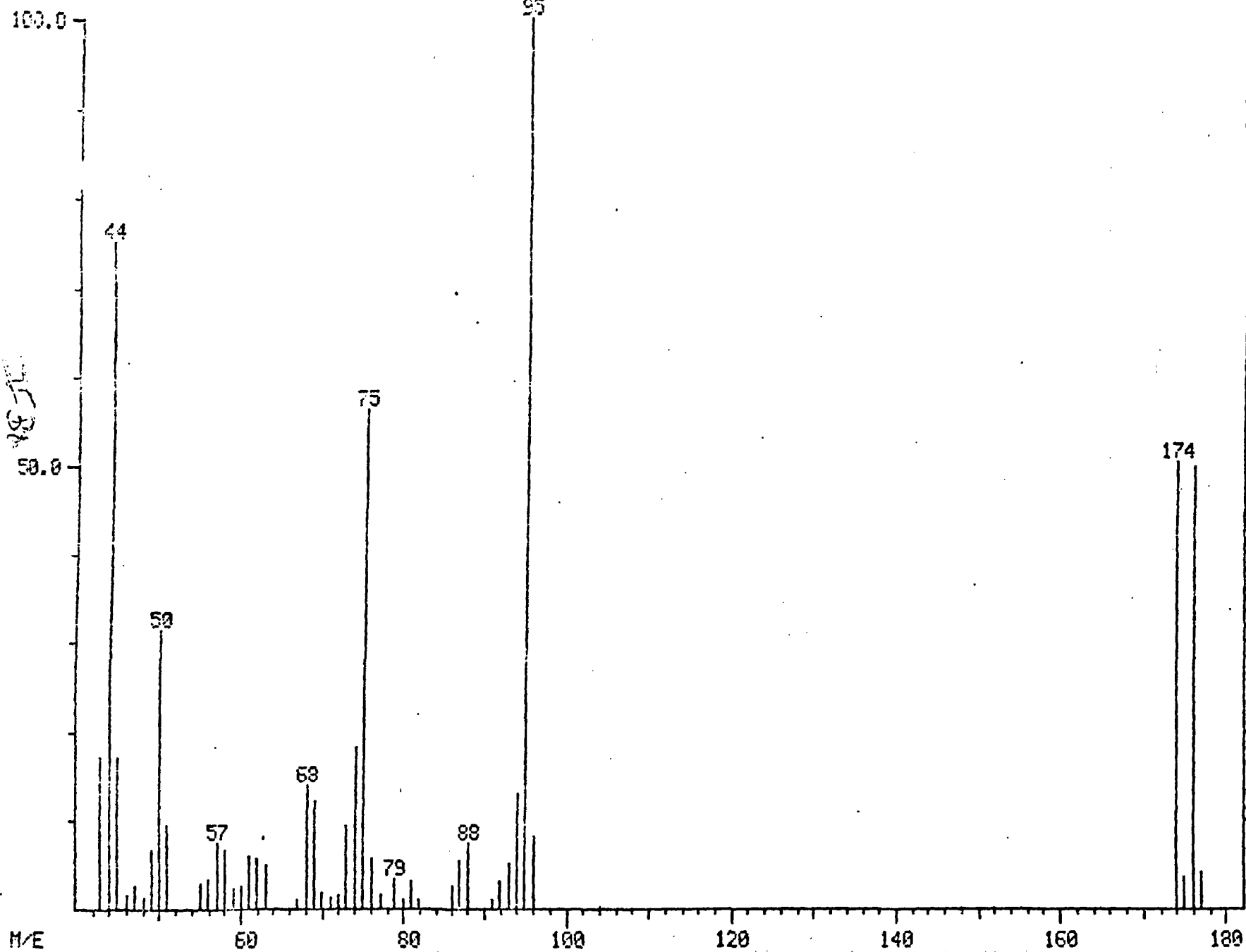
Mass	Rel. Abund.	Criterion
51	35.0688	30-60% MASS 198
68	0	< 2% MASS 69
69	66.2613	
70	.463632	< 2% MASS 69
127	58.3454	40-60% MASS 198
197	0	< 1% MASS 198
198	100	BASE PEAK
199	6.70366	5-9% MASS 198
275	17.5344	10-30% MASS 198
365	1.67971	> 1% MASS 198
441	0	< MASS 443
442	45.9755	> 40% MASS 198
443	8.34917	17-23% MASS 442

#-2

MASS SPECTRUM
03/04/05 6:41:00 + 30:16
SAMPLE: BLANK NBI CASE-9531-01-3769

DATA: V2715 #886

BASE M/E: 95
RIC: 62144.



10523.
10.

6:41:00 + 30:16
SAMPLE: BLANK MEL CASE ~~2715~~ ~~CM~~

3914
MIN INTEN: 0. MAX INTEN: 10529.

43 0.00 MINIMA
177 # 0 MAXIMA

MASS	%	RA
43	17.	08
44	75.	30
45	17.	10
46	1.	64
47	2.	58
48	1.	22
49	6.	70
50	21.	61
51	9.	43
55	2.	09
56	2.	33
57	7.	48
58	0.	33
59	2.	24
60	2.	55
61	5.	87
62	5.	64
63	4.	90
67	1.	13
68	14.	00
69	12.	37
70	1.	74
71	1.	40
72	1.	59
73	9.	48
74	18.	41
75	36.	31
76	5.	39
77	1.	36
79	3.	32
80	1.	01
81	3.	17
82	1.	03
86	2.	55
87	5.	33
88	7.	44
91	1.	06
92	2.	97
93	5.	01
94	12.	01
95	100.	00
96	8.	06
174	50.	30
175	3.	54
176	4.	85
177	4.	17

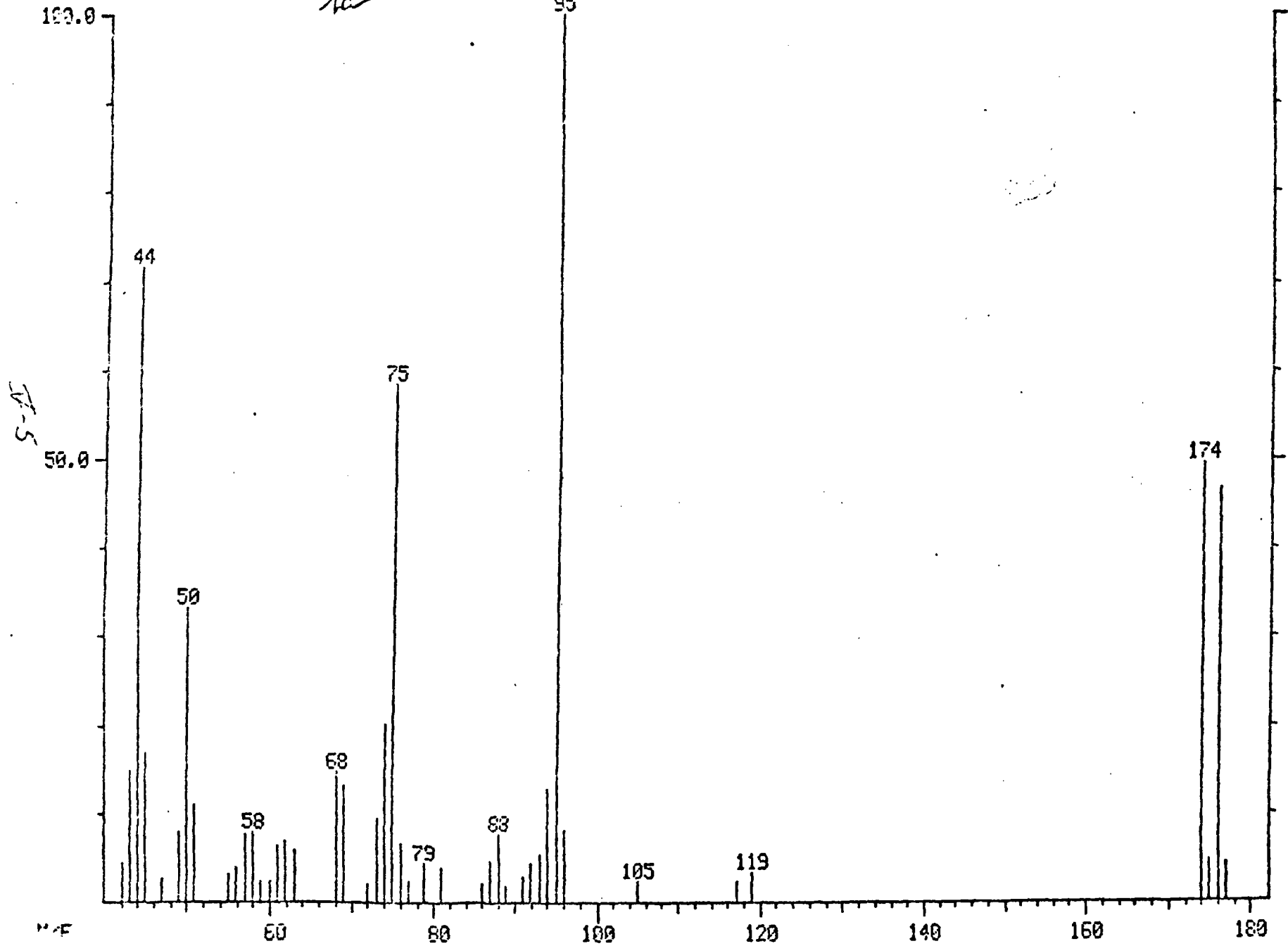
4

MASS SPECTRUM
03/05/85 17:40:00 + 30:20
SAMPLE: ~~7-3~~ 43236
ic

DATA: U2745 #888

BASE M/E: 95
RIC: 36932.

045 3969



5095.
10.

10/12

AMPLE: 43236

Case 3769

42 0.00 MINIMA MIN INTEN: 0. MAX INTEN: 6096.
 177 0 MAXIMA
 MASS % RA

42	4.22
43	14.78
44	71.65
45	16.93
47	2.44
49	7.94
50	33.30
51	10.94
55	3.15
56	3.89
57	7.68
58	7.86
59	2.18
60	2.31
61	6.40
62	6.96
63	5.87
68	14.80
69	13.39
72	2.00
73	9.40
74	20.24
75	33.53
76	6.66
77	2.30
79	4.27
81	3.77
86	2.12
87	4.64
88	7.73
89	1.74
91	2.77
92	4.23
93	5.33
94	17.88
95	100.00
96	8.19
105	2.40
117	2.30
119	3.25
174	49.61
175	4.66
176	46.72
177	4.22

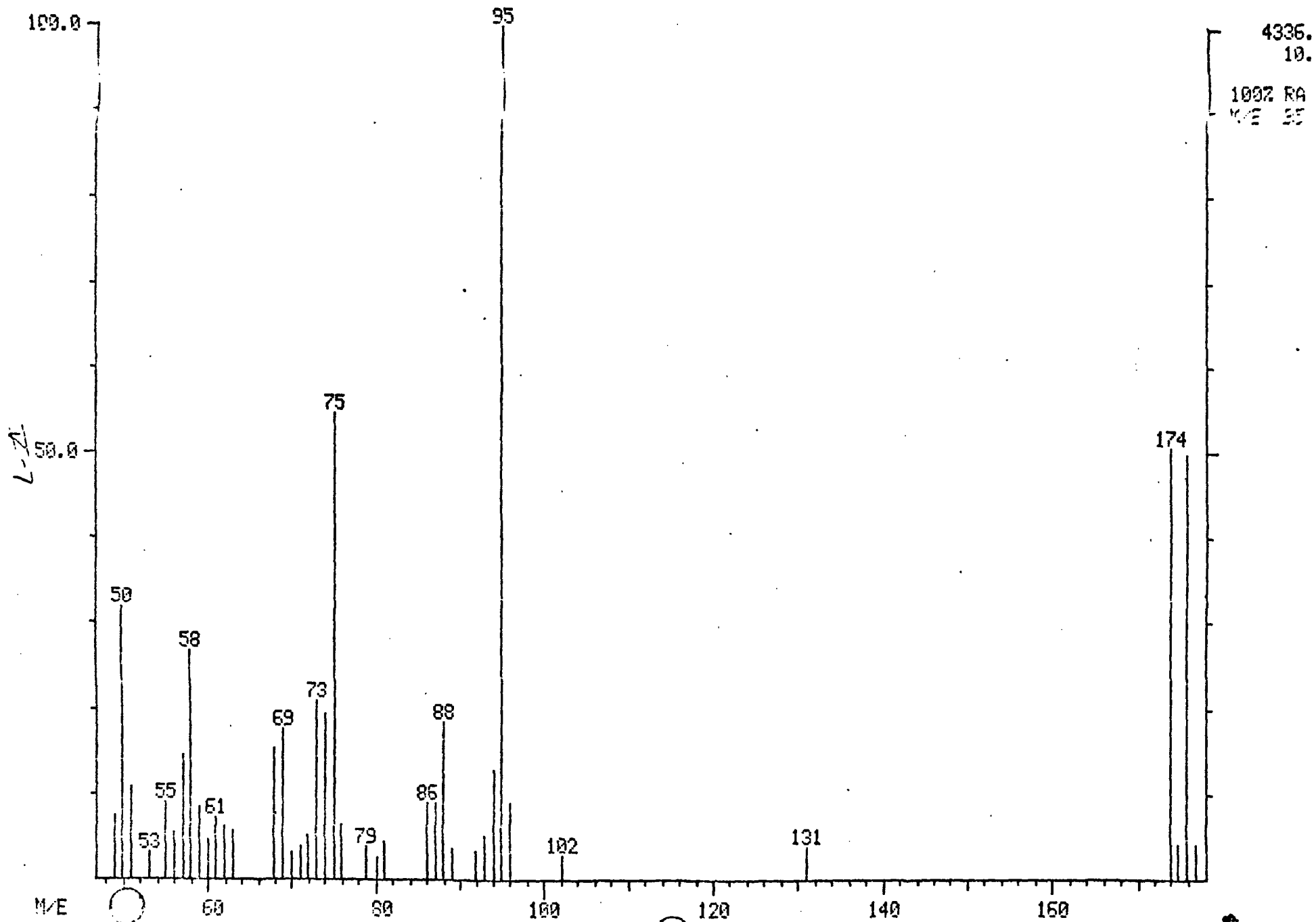
P-6

20

MASS SPECTRUM
03/05/85 10:26:00 + 9:57
SAMPLE: BFB 3/5/85

DATA: U2753 #291

BASE M/E: 44
RIC: 41024.



4336.
10.
100% RA
M/E 35

ASS	#	% RA	MINIMA	MAXIMA
48	0.00		MINIMA	
177	0		MAXIMA	
49	7.36			
50	31.83			
51	10.77			
53	3.02			
55	8.86			
56	5.33			
57	14.48			
58	26.71			
59	8.46			
60	4.70			
61	7.06			
62	6.04			
63	5.58			
68	15.29			
69	17.53			
70	3.16			
71	3.94			
72	5.17			
73	20.85			
74	19.44			
75	54.61			
76	6.37			
79	3.76			
80	2.68			
81	4.29			
86	8.97			
87	8.93			
88	18.27			
89	3.69			
92	3.41			
93	5.14			
94	12.66			
95	100.00			
96	9.04			
102	2.70			
131	3.85			
174	50.55			
175	4.13			
176	47.63			
177	4.06			

MIN INTEN: 0. MAX INTEN: 4336.
100% RA AT M/E: 95

V-8

22

Laboratory Name: GCA
 Lab Sample ID No: Blank
 Sample Matrix: Water
 Data Release Authorized By: Fiona Marquis

Case No. 3931
 QC Report No. 3931
 Contract No. 68-01-1767
 Date Sample Received: 3/4/85

Volatile Compounds

Concentration Low Medium (Circle One)
 Date Extracted/Prepared: 3/4/85
 Date Analyzed: 3/4/85
 Conc/Dil Factor: 1 pH: NA
 Percent Moisture: NA
 Percent Moisture (Decanted): NA

CAS Number	Compound	ug/l or ug/Kg (Circle One)
74-87-3	Chloroethane	10u
74-83-8	Bromomethane	10u
75-01-4	Vinyl Chloride	10u
75-00-3	Chloroethane	10u
75-10-2	Methylene Chloride	5u
67-54-1	Acetone	10u
75-15-0	Carbon Disulfide	5u
75-35-4	1,1-Dichloroethane	5u
75-24-3	1,1-Dichloroethane	5u
156-09-5	Trans-1,2-Dichloroethane	5u
67-16-3	Chloroform	5u
107-06-2	1,2-Dichloroethane	5u
78-53-3	1-Butanone	10u
71-55-6	1,1,1-Trichloroethane	5u
11-23-5	Carbon Tetrachloride	5u
103-05-4	Vinyl Acetate	10u
75-27-4	Bromodichloromethane	5u

CAS Number	Compound	ug/l or ug/Kg (Circle One)
73-34-6	1,1,2,2-Tetrachloroethane	5u
73-87-5	1,2-Dichloropropane	5u
10061-02-6	Trans-1,3-Dichloropropene	5u
79-01-6	Trichloroethane	5u
124-48-1	Dibromochloromethane	5u
79-00-5	1,1,2-Trichloroethane	5u
71-43-2	Benzene	5u
10461-01-5	cis-1,3-Dichloropropene	5u
110-75-0	2-Chloroethylvinylether	10u
75-25-2	Bromotoluene	5u
591-78-6	2-Hexanone	10u
108-10-1	4-Methyl-2-Pentanone	10u
127-18-4	Tetrachloroethene	5u
109-88-3	Toluene	5u
109-90-7	Chlorobenzene	5u
100-41-4	Ethylbenzene	5u
100-42-5	Styrene	5u
	Total Xylenes	5u

Environmental Qualifiers

For a complete explanation of the following qualifiers, consult the manual.
 Additional help is available by explaining the environmental. However, the
 definition of each qualifier is as follows:

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

U: Indicates concentration was less than the detection limit. Report the minimum detectable level for the sample (e.g., 100) based on necessary concentration factor among (U) is not necessarily the minimum detection limit. The detection limit is (U). Compound was analyzed for that sample but the number in the next column is the detection limit for the sample.

J: Indicates an estimated value. This flag is used when the concentration is estimated to be statistically significant compared to when the response is zero or when the response is not statistically significant. The number in the next column is the detection limit for the sample.

C: This flag is used to indicate parameters where the detection limit has been confirmed by GC/MS single component detection. All ug/l or the four digit standard deviation by GC/MS.

P: This flag is used to indicate analysis of a sample in the field as well as the sample indicates possible possible detection limit. It warns the detector to take appropriate action.

D: This flag is used to indicate parameters where the detection limit has been confirmed by GC/MS single component detection. All ug/l or the four digit standard deviation by GC/MS.

see manual

IV-9

Organics Analysis Data Sheet
(Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 3/5/85

Date Analyzed: 3/25/85

Conc/Dil Factor: 1.0L → 10mL

CAS Number		<u>ug/l</u> or ug/Kg (Circle One)
62-75-9	N-Nitrosodimethylamine	10u
108-95-2	Phenol	10u
62-53-3	Aniline	10u
111-44-4	Di-2-Chloroethyl Ether	10u
95-57-8	2-Chlorophenol	10u
541-73-1	1,3-Dichlorobenzene	10u
106-46-7	1,4-Dichlorobenzene	10u
100-51-6	Benzyl Alcohol	10u
95-50-1	1,2-Dichlorobenzene	10u
95-48-7	2-Methylphenol	10u
39638-32-9	Di-2-chloroisopropyl ether	10u
106-44-5	4-Methylphenol	10u
621-64-7	N-Nitroso-Di-n-Propylamine	10u
67-72-1	Hexachlorocyclopentadiene	10u
98-95-3	Nitrobenzene	10u
78-59-1	Heptachloroepoxide	10u
88-75-5	2-Nitrophenol	10u
105-67-9	2,4-Dimethylphenol	10u
65-85-0	Benzoic Acid	50u
111-91-1	Di-2-Chloro-isopropyl Methane	10u
120-83-2	2,4-Dichlorophenol	10u
120-82-1	1,2,4-Trichlorobenzene	10u
91-20-3	Fluoranthene	10u
106-47-6	4-Chloroaniline	10u
87-68-3	Hexachlorobenzene	10u
59-50-7	4-Chloro-3-Methylphenol	10u
91-57-6	2-Methylnaphthalene	10u
77-47-4	Hexachlorocyclopentadiene	10u
88-06-2	2,4,6-Trichlorophenol	10u
95-95-4	2,4,5-Trichlorophenol	10u
91-58-7	2-Chloronaphthalene	10u
88-74-4	2-Nitroaniline	10u
131-11-3	Dimethyl Phthalate	10u
203-96-6	Acenaphthylene	10u
98-09-2	5-Nitroaniline	10u

CAS Number		<u>ug/l</u> or ug/Kg (Circle One)
33-32-9	Acenaphthone	10u
51-28-5	2,4-Dinitrophenol	10u
100-02-7	4-Nitrophenol	10u
132-64-9	Dibenzofuran	10u
121-14-2	2,4-Dinitrotoluene	10u
606-20-2	2,6-Dinitrotoluene	10u
84-66-2	Diethylphthalate	10u
7005-72-3	4-Chlorobenzyl phenylether	10u
95-73-7	Fluorene	10u
100-01-6	4-Nitroaniline	10u
534-52-1	4,6-Dinitro-2-Methylphenol	10u
26-30-6	N-Nitrosodiphenylamine (1)	10u
101-55-3	4-Bromophenyl phenylether	10u
118-74-1	Hexachlorobenzene	10u
37-85-5	Pentachlorophenol	10u
85-01-8	Phenanthrene	10u
120-12-7	Anthracene	10u
34-74-2	Di-n-Butylphthalate	10u
206-44-0	Fluoranthene	10u
92-87-5	Benzidine	10u
129-00-0	Pyrene	10u
85-68-7	Dibutylziphthalate	10u
91-94-1	3,3'-Dichlorobenzidine	10u
96-55-3	Benzofluoranthene	10u
117-81-7	Di-2-Ethylhexyl Phthalate	10u
118-01-9	Chrysene	10u
117-84-0	Di-n-Octyl Phthalate	10u
205-99-2	Benzofluoranthene	10u
207-08-9	Benzokluoranthene	10u
50-32-8	Benzo(a)Pyrene	10u
193-50-5	Indeno(1,2,3-cd)pyrene	10u
53-70-3	Dibenz(a,h)Anthracene	10u
191-24-2	Benzo(g,h,i)perylene	10u

(1) Cannot be separated from diphenylamine

Sample Number
Method Blank

Organics Analysis Data Sheet
 (Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 3-1-85

Date Analyzed: 3/30/85

Conc/Dil Factor: 1

Michael 4/13/85

CAS Number ug/l or ug/Kg
(Circle One)

319-84-6	Alpha-BHC	0.05U
319-85-7	Beta-BHC	0.05U
319-86-8	Delta-BHC	0.05U
58-89-9	Gamma-BHC (Lindane)	0.05U
76-44-8	Heptachlor	0.05U
309-00-2	Aldrin	0.05U
1024-57-3	Heptachlor Epoxide	0.05U
959-98-8	Endosulfan I	0.05U
60-57-1	Dieldrin	0.10U
72-55-9	4, 4'-DDE	0.10U
72-20-8	Endrin	0.10U
33213-65-9	Endosulfan II	0.10U
72-54-8	4, 4'-DDD	0.10U
7421-93-4	Endrin Aldehyde	0.10U
1031-07-8	Endosulfan Sulfate	0.10U
50-29-3	4, 4'-DDT	0.10U
72-43-6	Methoxychlor	0.50U
53494-70-5	Endrin Ketone	0.10U
57-74-9	Chlordane	0.50U
8001-35-2	Toxaphene	1.0U
12674-11-2	Aroclor-1615	0.50U
11104-28-2	Aroclor-1231	0.50U
11141-16-5	Aroclor-1232	0.50U
53469-21-9	Aroclor-1242	0.50U
12672-29-6	Aroclor-1248	0.50U
11097-69-1	Aroclor-1254	1.0U
11096-82-5	Aroclor-1260	1.0U



V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s 1000 ml or W_s _____ V_i 10000 ul V_t 5.0 ul

Organics Analysis Data Sheet (Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	<i>No compounds found</i>	<i>VQA</i>		
2.				
3.				
4.				
5.				
6.				
6.				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
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21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				

10-12-201

[Handwritten signature]

Organics Analysis Data Sheet
(Page 4)

MB1

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	2,4-dinitrophenol	BX1A	24.9	21 J
2.				
3.				
4.				
5.				
6.				
7.				
8.				
9.				
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48.				
49.				
50.				

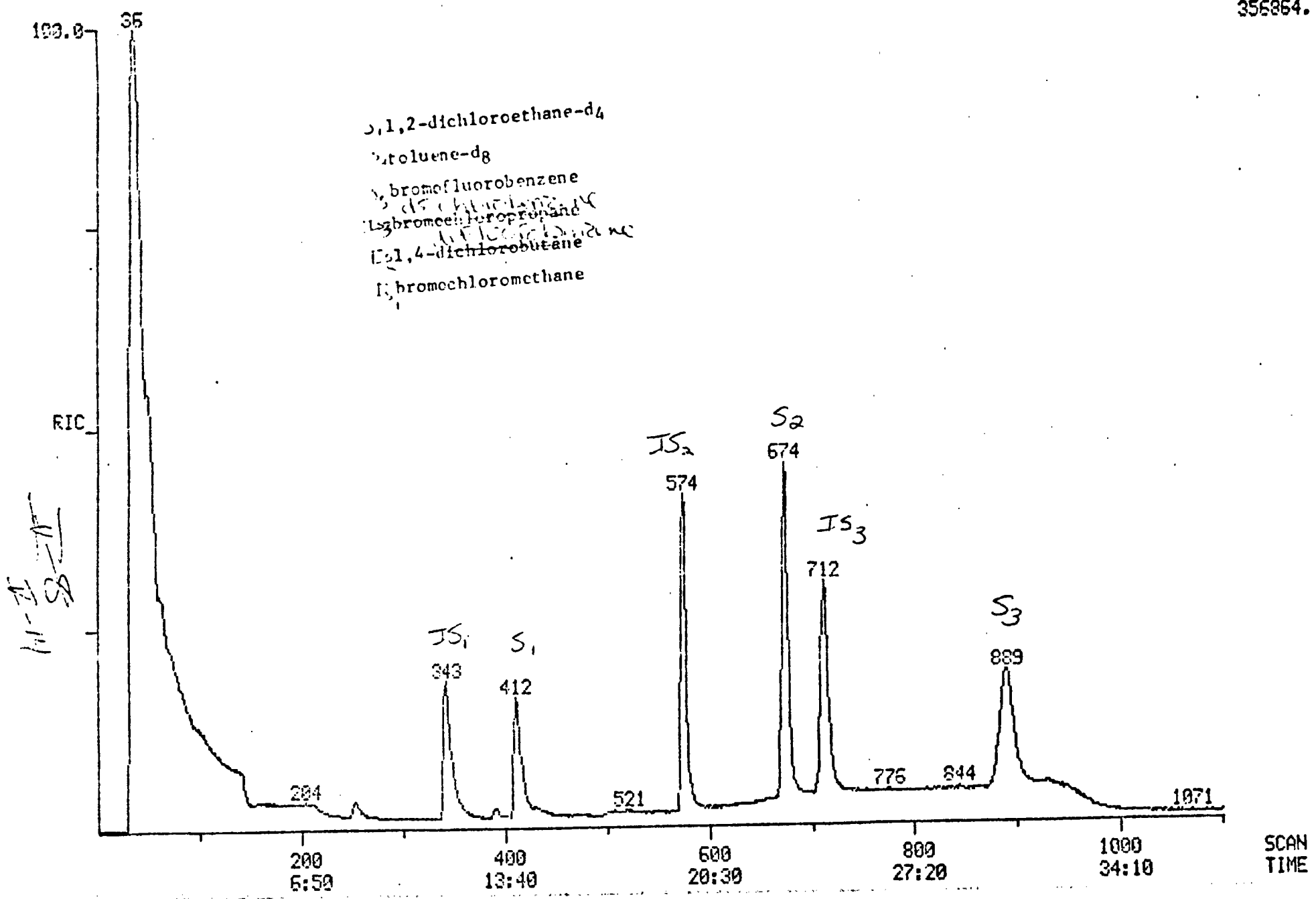
Page 4 of 4 11-29 II-13 11/89

RIC
03/04/85 6:41:00
SAMPLE: BLANK MS1 CASE 3931 CM

DATA: U2715

SCANS 1 TO 1100

356864.



Case:

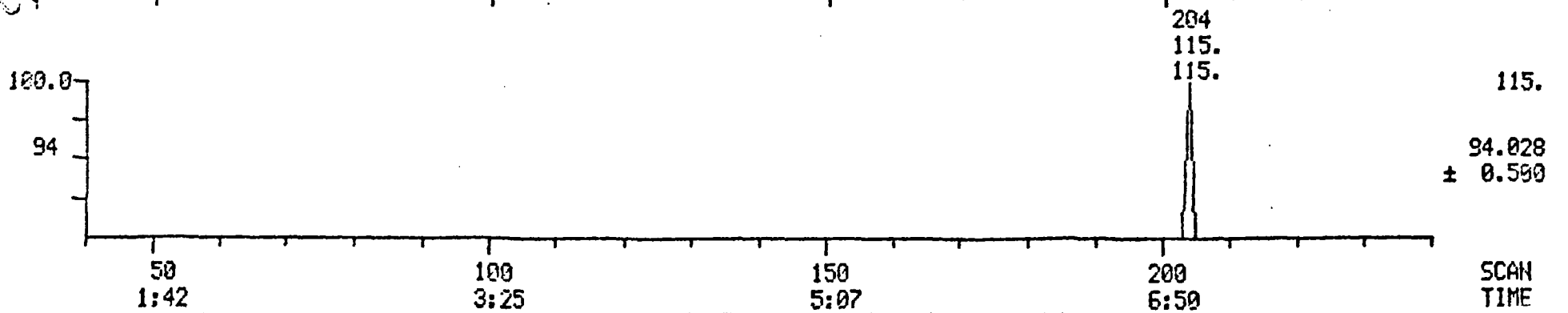
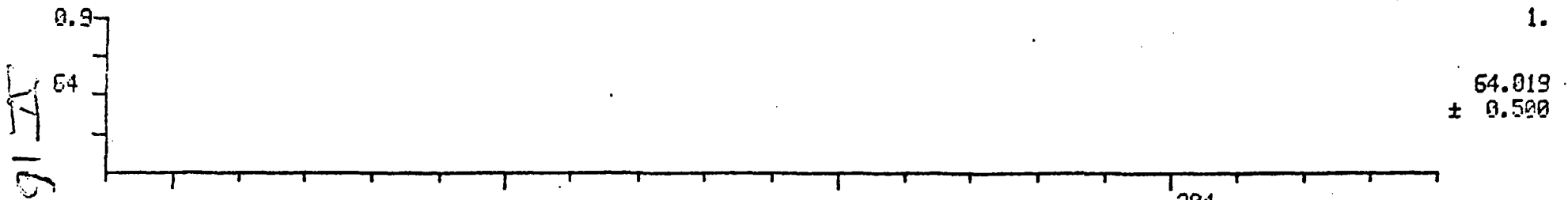
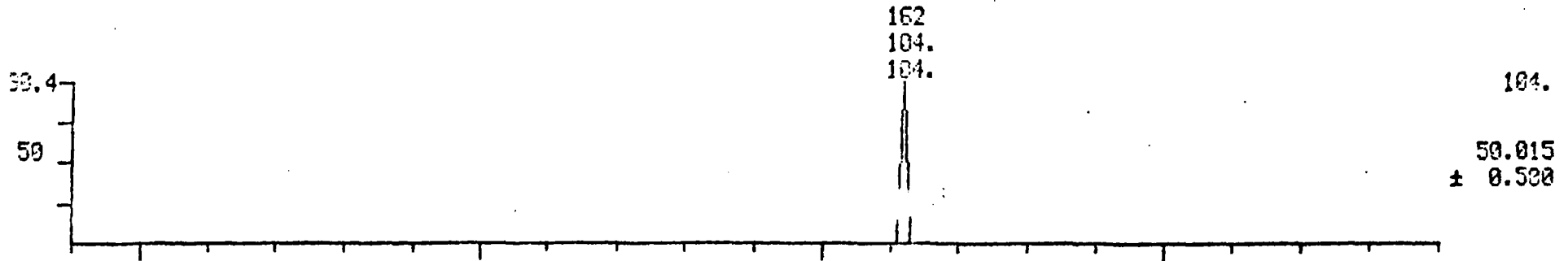
3969

NO	NAME
1	BROMOCHLOROMETHANE (I. S. #1)
2	CHLOROMETHANE
3	BROMOMETHANE
4	VINYL CHLORIDE
5	CHLOROETHANE
6	METHYLENE CHLORIDE
7	CARBON DISULFIDE
8	ACETONE
9	1,1-DICHLOROETHENE
10	1,1-DICHLOROETHANE
11	TRANS-1,2-DICHLOROETHENE
12	CHLOROFORM
13	D4-1,2-DICHLOROETHANE (SURR #1)
14	1,2-DICHLOROETHANE
15	1,1,1-TRICHLOROETHANE
16	CARBON TETRACHLORIDE
17	VINYL ACETATE
18	2-BUTANONE
19	1,4-difluorobenzene (I. S. #2)
20	BROMODICHLOROMETHANE
21	1,2-DICHLOROPROPANE
22	TRANS-1,3-DICHLOROPROPENE
23	TRICHLOROETHENE
24	DIBROMOCHLOROMETHANE
25	1,1,2-TRICHLOROETHANE
26	CIS-1,2-DICHLOROPROPENE
27	BENZENE
28	2-CHLOROETHYL VINYL ETHER
29	BROMOFORM
30	D8-TOLUENE
31	D5-CHLOROBENZENE (I. S. #3)
32	TETRACHLOROETHENE
33	1,1,2,2-TETRACHLOROETHANE
34	2-METHYL-2-PENTANONE
35	2-HEXANONE
36	TOLUENE
37	CHLOROBENZENE
38	ETHYL BENZENE
39	BFB (SURR #3)
40	STRYENE
41	TOTAL XYLENES

MASS CHROMATOGRAMS
03/04/85 5:41:00
SAMPLE: BLANK NB1 CASE 3331 CM

DATA: U2715

SCANS 40 TO 240



50
1:42

100
3:25

150
5:07

200
6:50

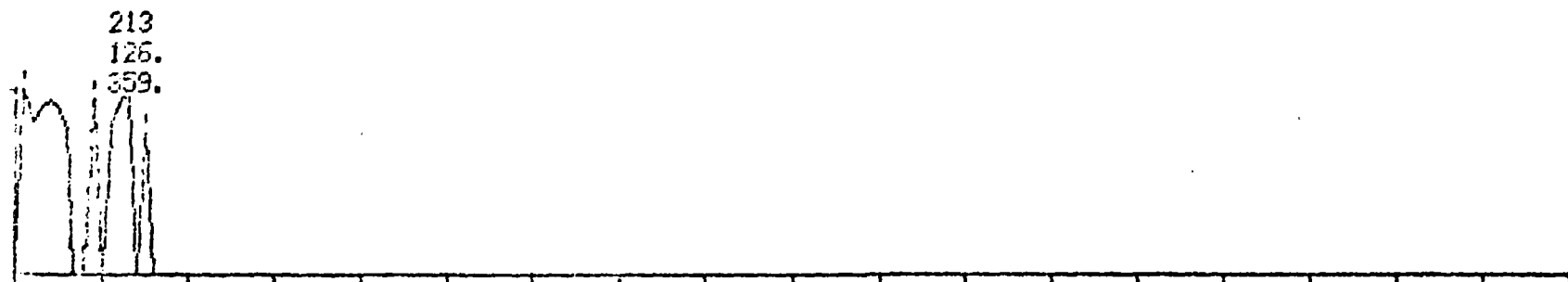
SCAN
TIME

MASS CHROMATOGRAMS
23/04/85 6:41:00
SAMPLE: BLANK NDI CASE 3331 CM

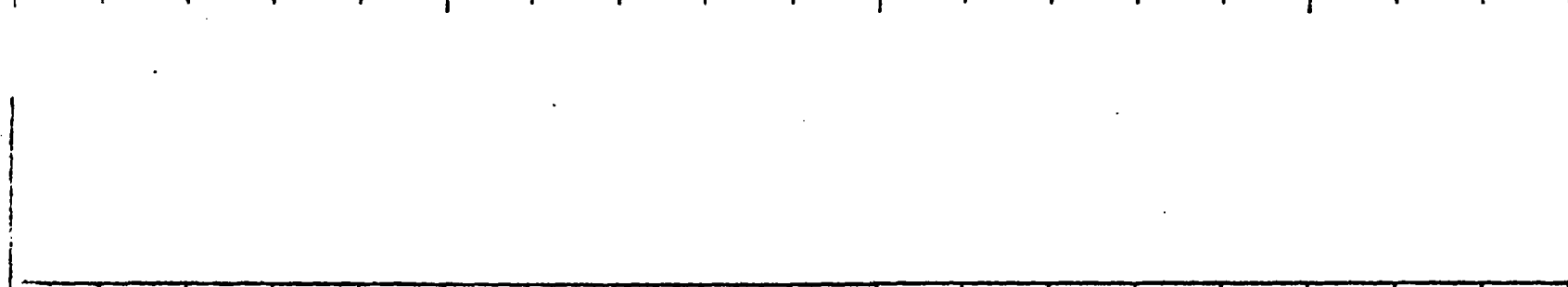
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SCANS 200 TO 330

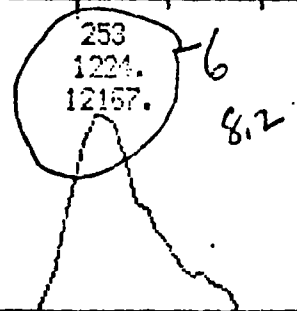
U-7



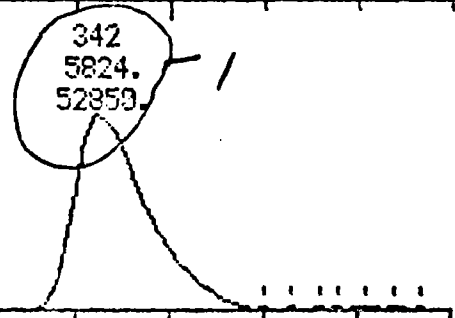
131.
58.917
± 0.500



1.
76.023
± 0.500



1224.
84.025
± 0.500



5824.
128.038
± 0.500

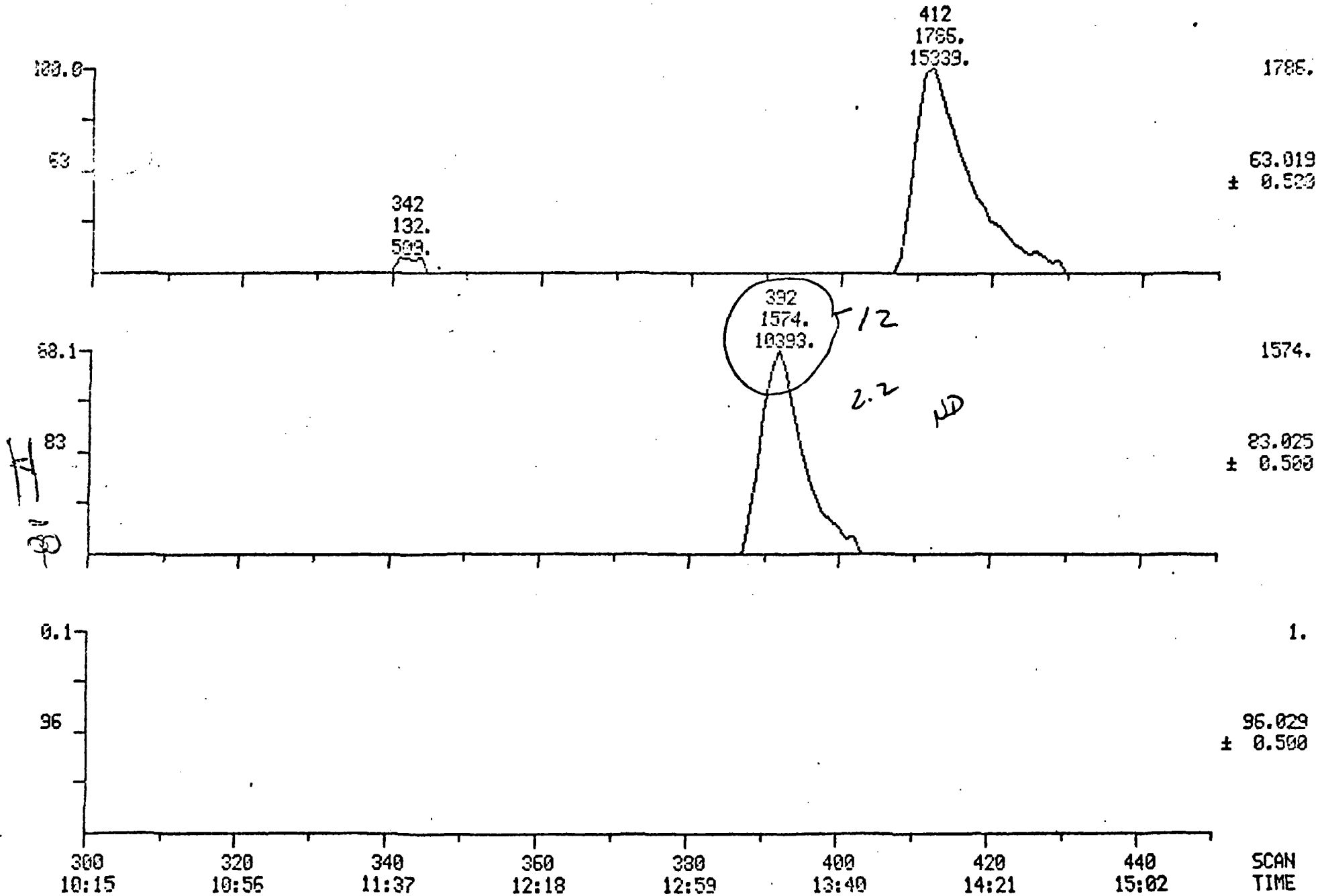
250 8:32 300 10:15 350 11:57

SCAN
TIME

MASS CHROMATOGRAMS
03/04/85 6:41:00
SAMPLE: BLANK 181 CASE 3931 CM

DATA: U2715

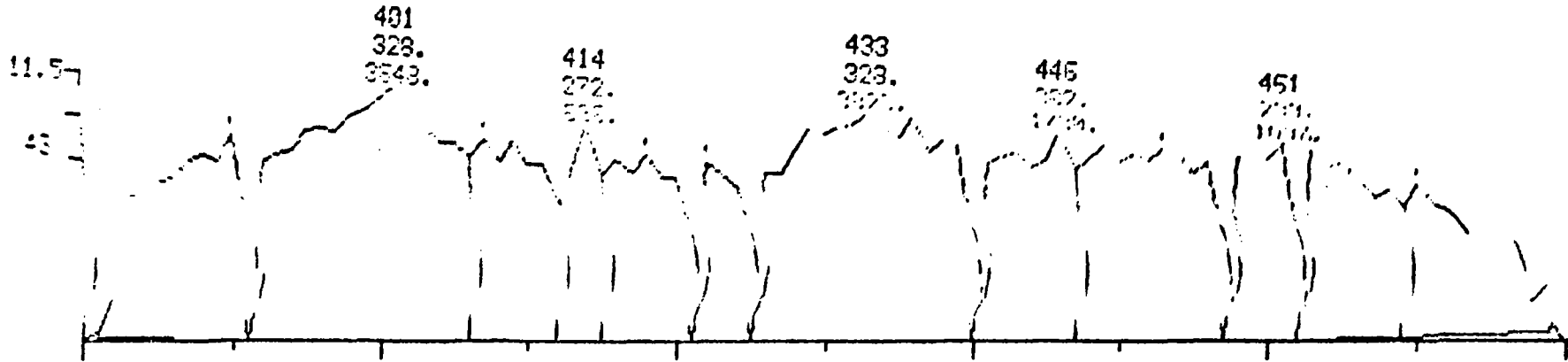
SCANS 389 TO 450



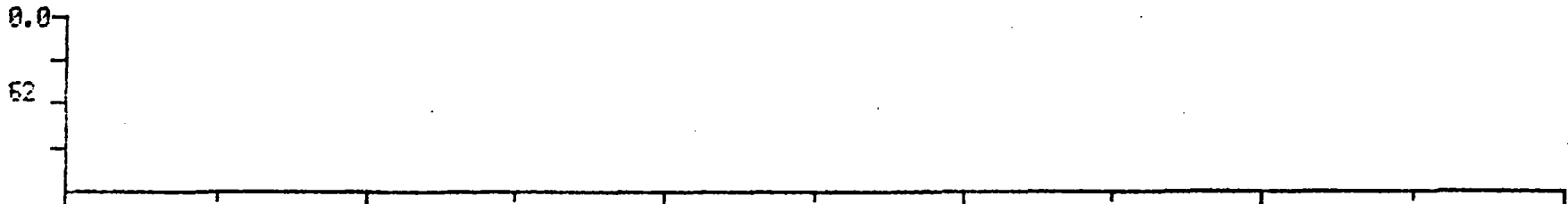
55 CHROMATOGRAMS
03/04/85 6:41:00
SAMPLE: BLANK NB1 CASE 3931 CM

DATA: U2715

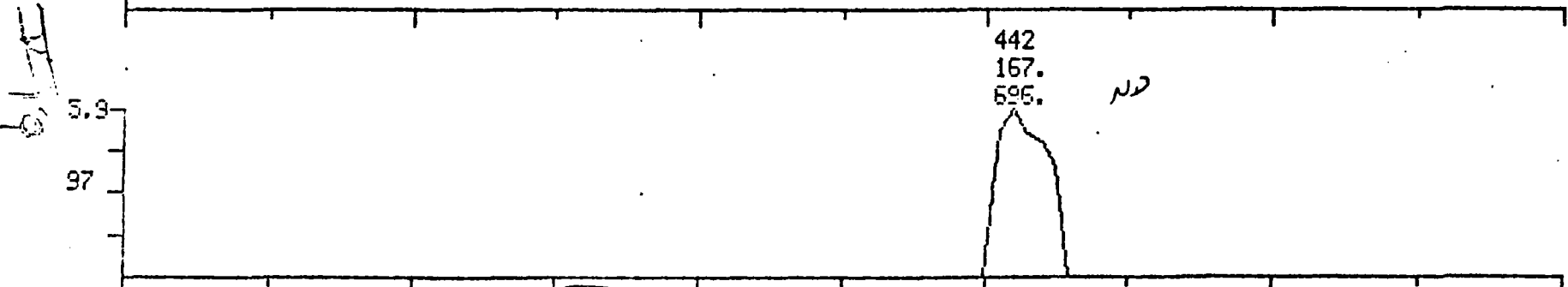
SCANS 390 TO 480



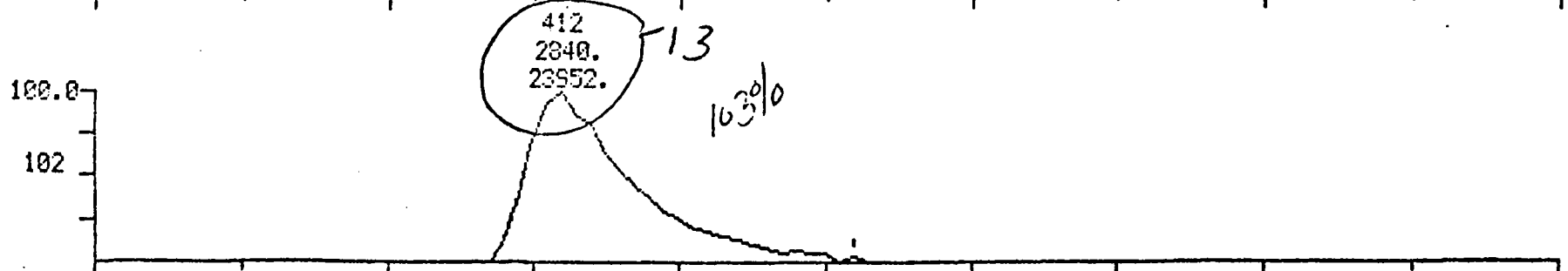
40.213
± 0.500



1.
52.019
± 0.500



157.
97.029
± 0.500



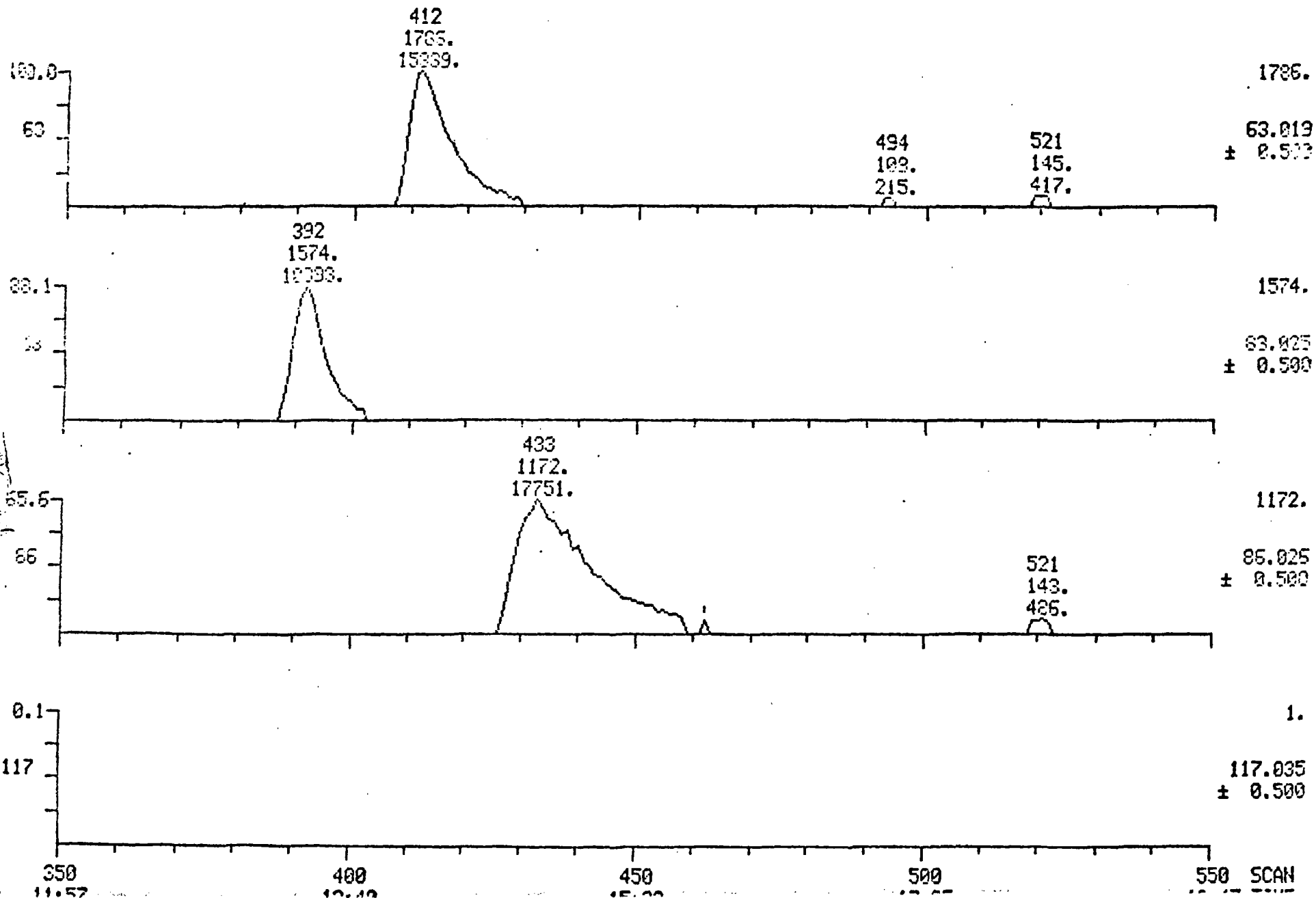
2840.
102.031
± 0.500

390 12:59 400 13:40 420 14:21 440 15:02 460 15:43 480 15:24 SCAN TIME

MASS CHROMATOGRAMS
03/04/85 6:41:20
SAMPLE: BLANK MB1 CASE 3931 CM

DATA: U2715

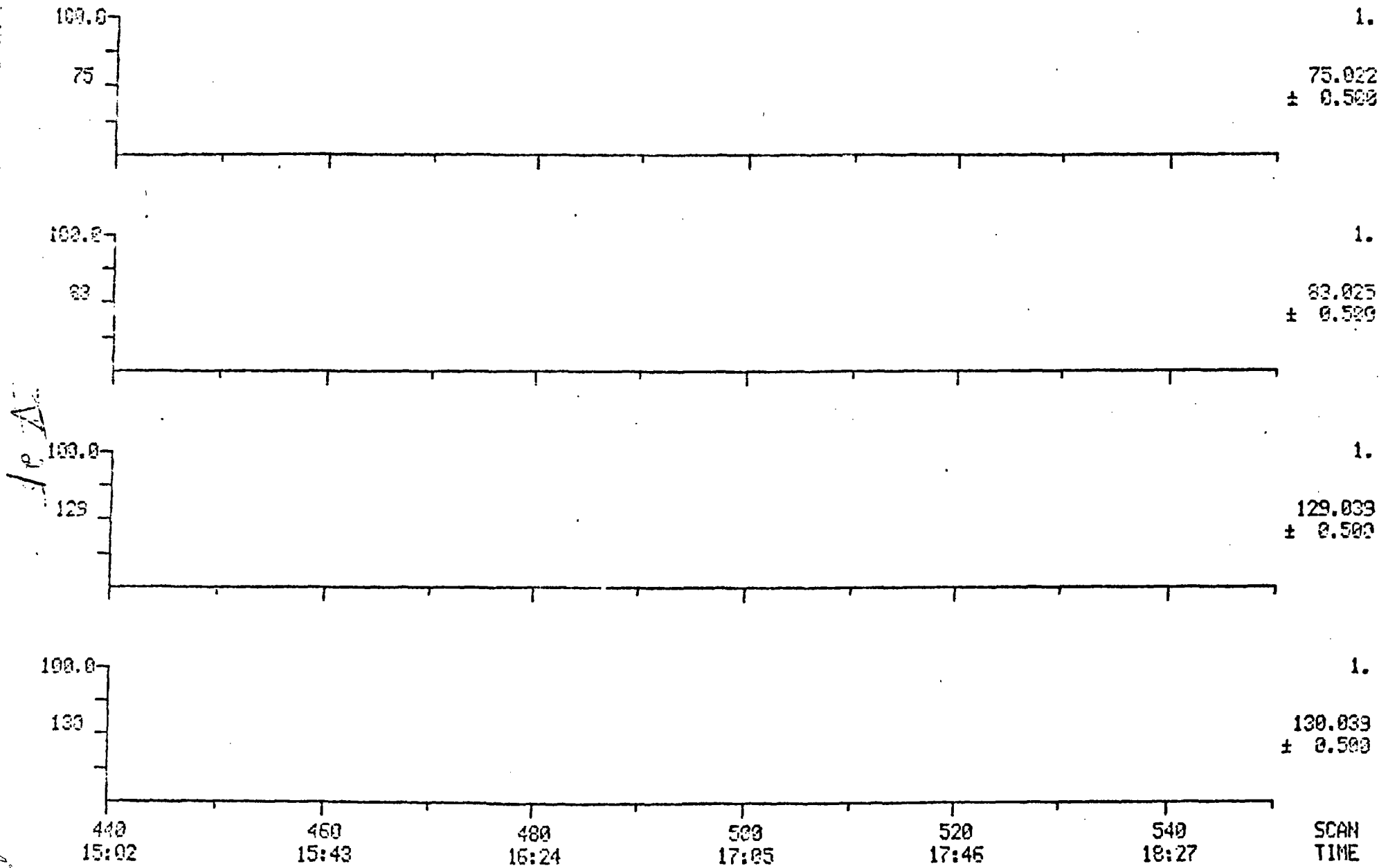
SCANS 350 TO 550



MASS CHROMATOGRAMS
03/04/05 6:41:00
SAMPLE: BLANK FBI CASE 3931 CM

DATA: U2715

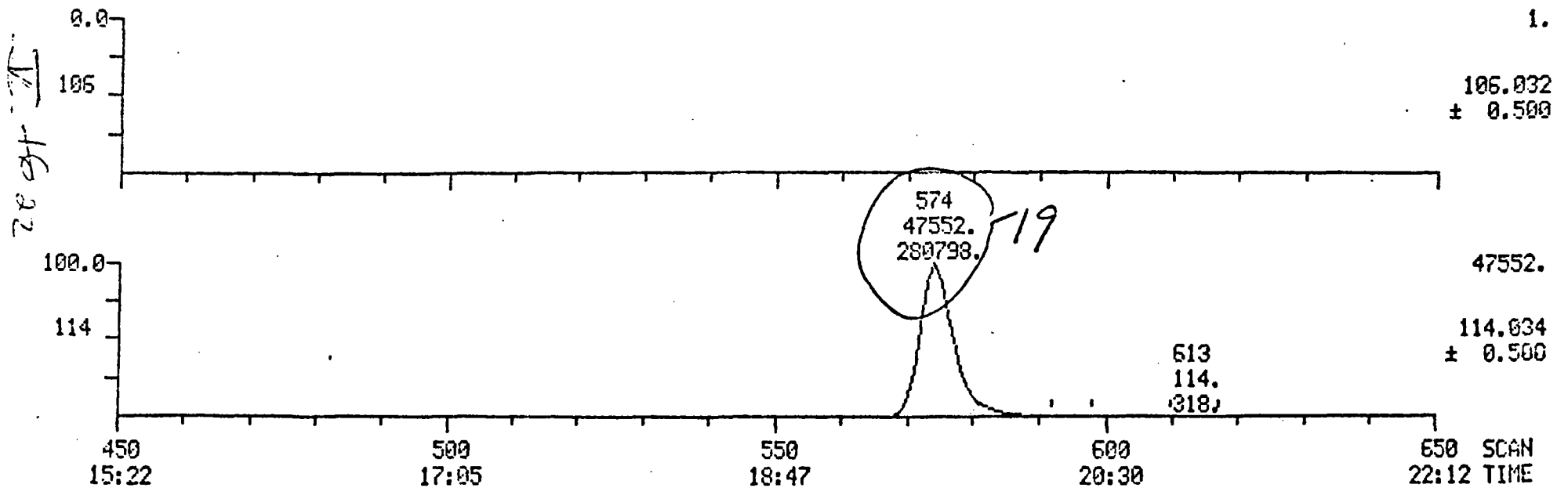
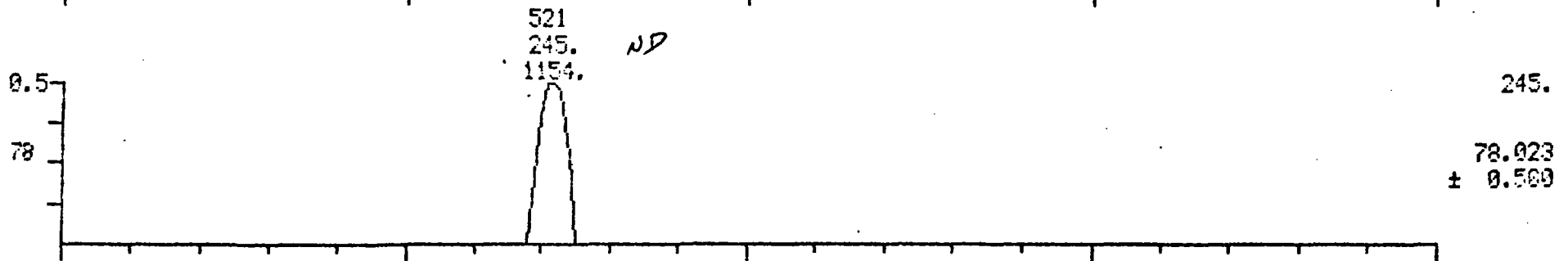
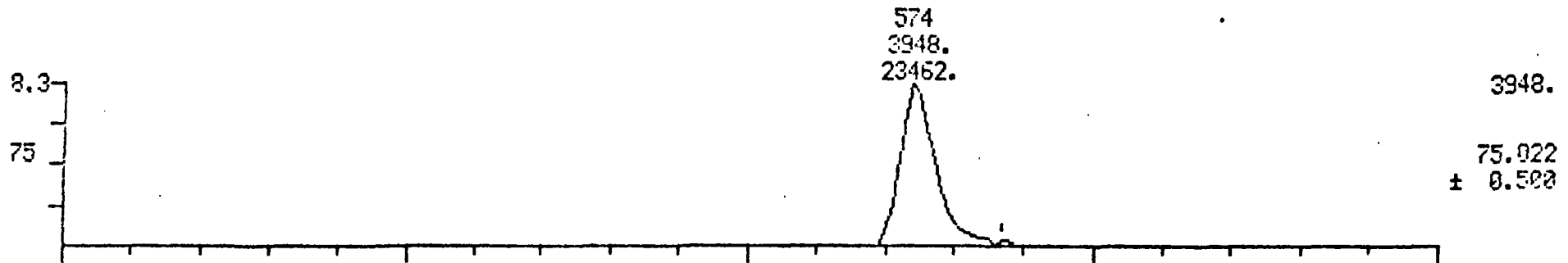
SCANS 440 TO 550



MASS CHROMATOGRAMS
03/04/85 6:41:00
SAMPLE: BLANK NBI CASE 3931 CM

DATA: U2715

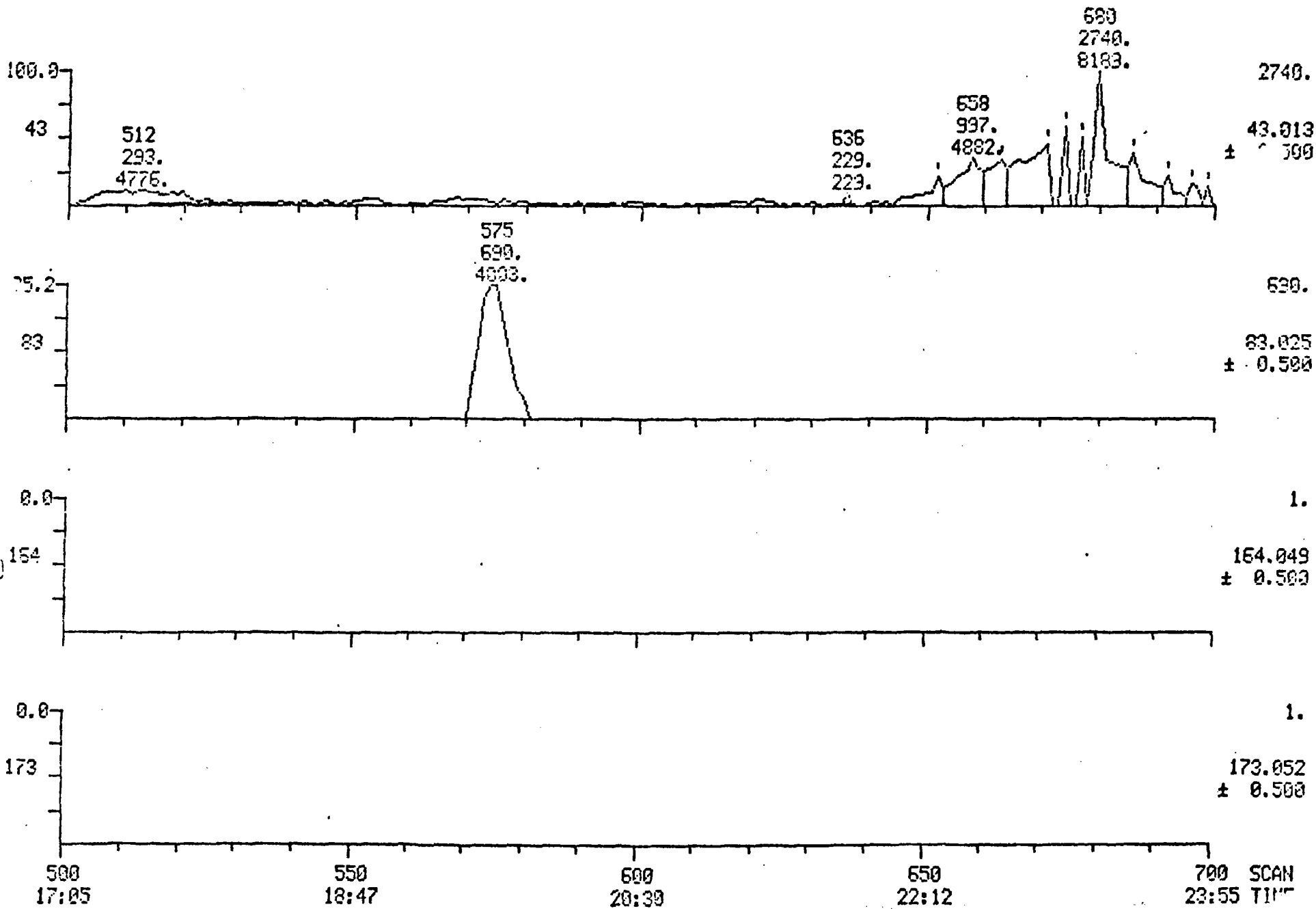
SCANS 450 TO 650



MASS CHROMATOGRAMS
03/04/05 6:41:00
SAMPLE: BLANK MBI CASE 3931 CM

DATA: U2715

SCANS 500 TO 700



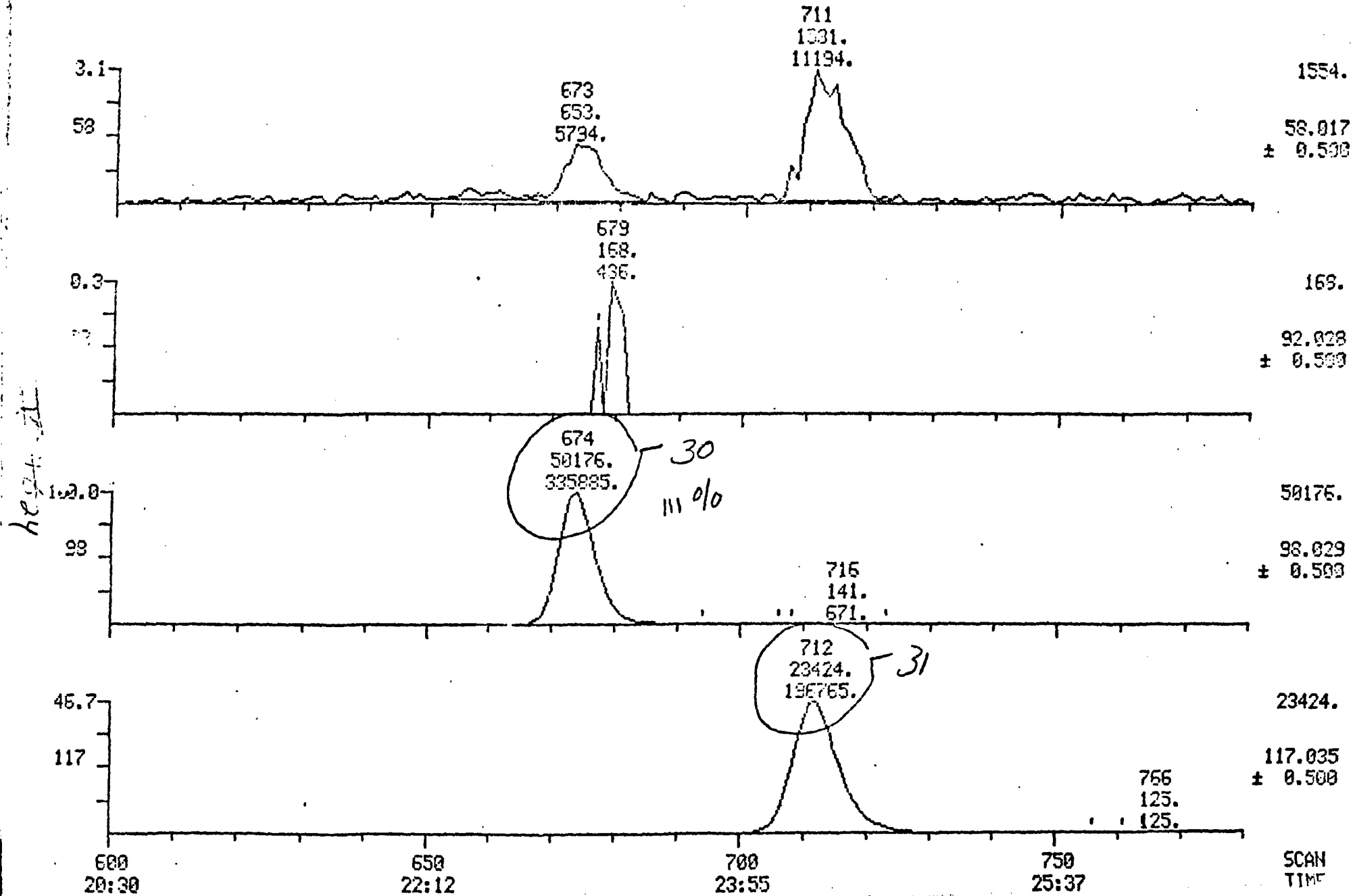
3071-1703

23

MASS CHROMATOGRAMS
03/04/85 6:41:00
SAMPLE: BLANK NB1 CASE 3931 CM

DATA: U27.5

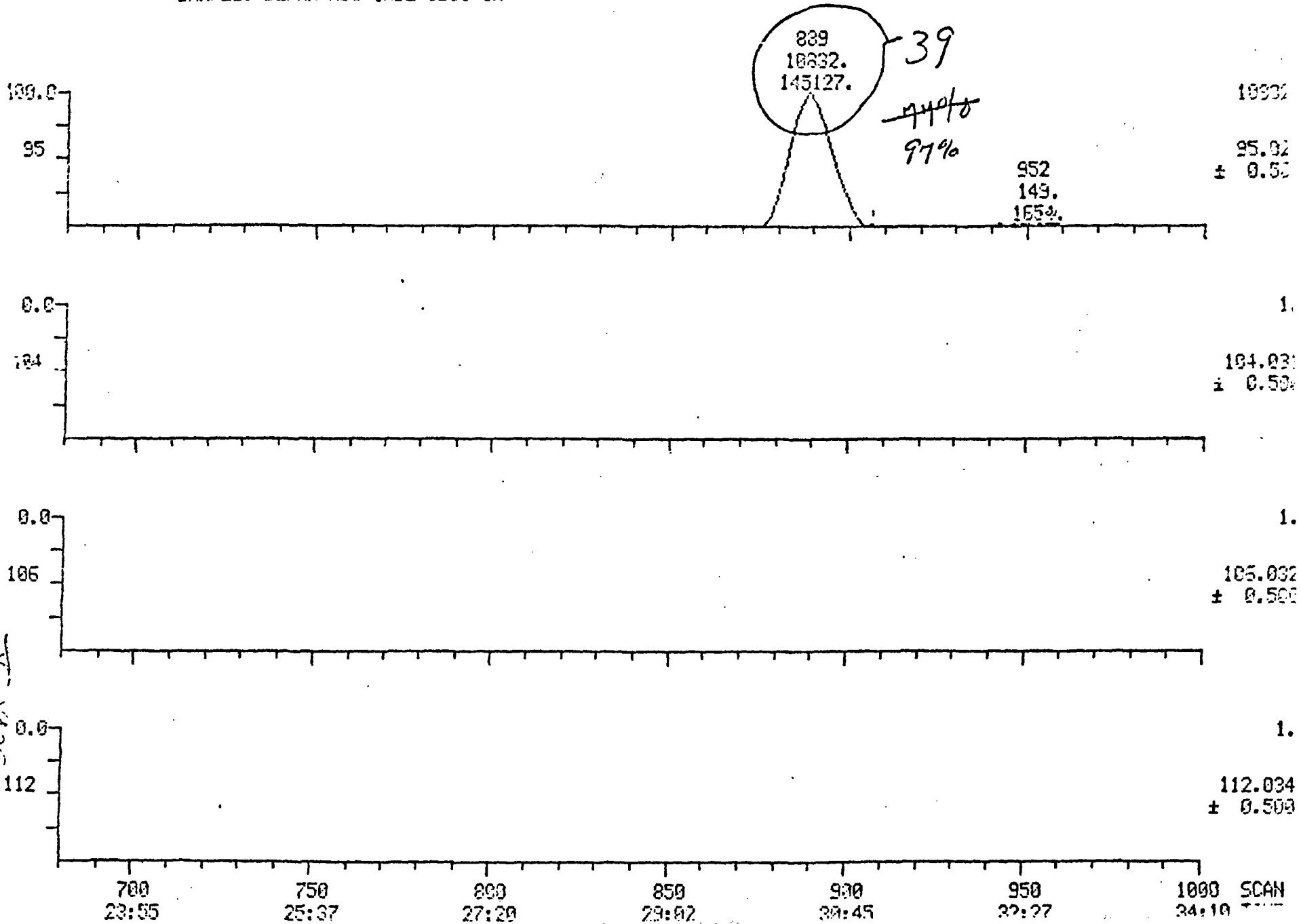
SCANS 600 TO 720



03/04/85 6:41:00
SAMPLE: BLANK ME1 CASE 3931 CM

DATA: 01715

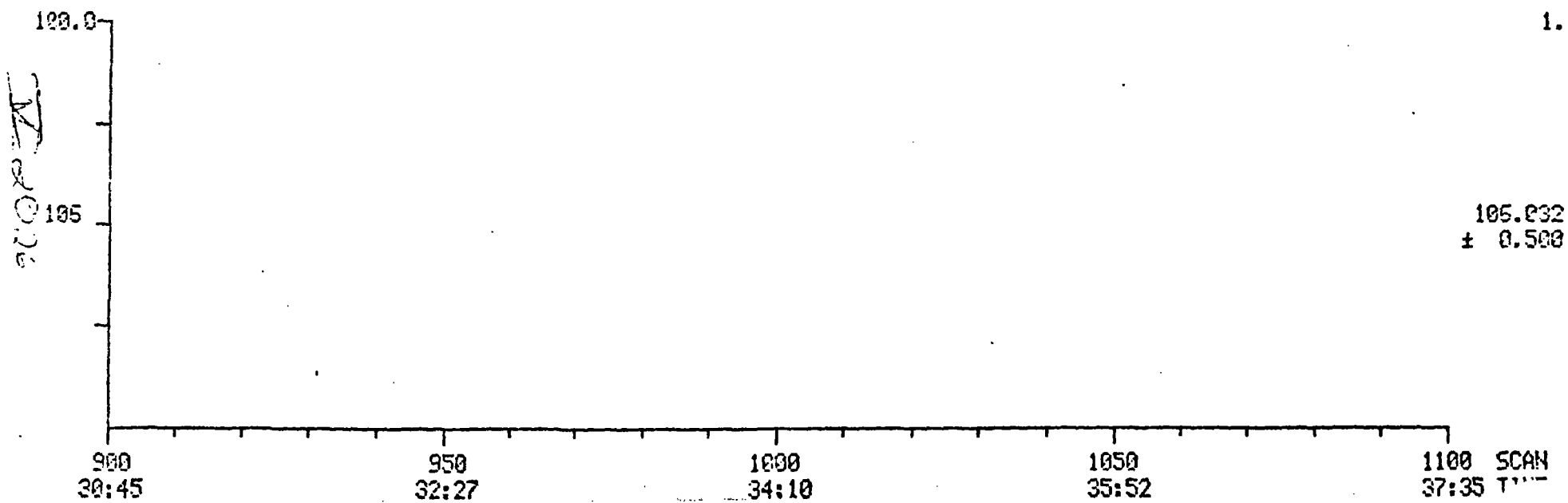
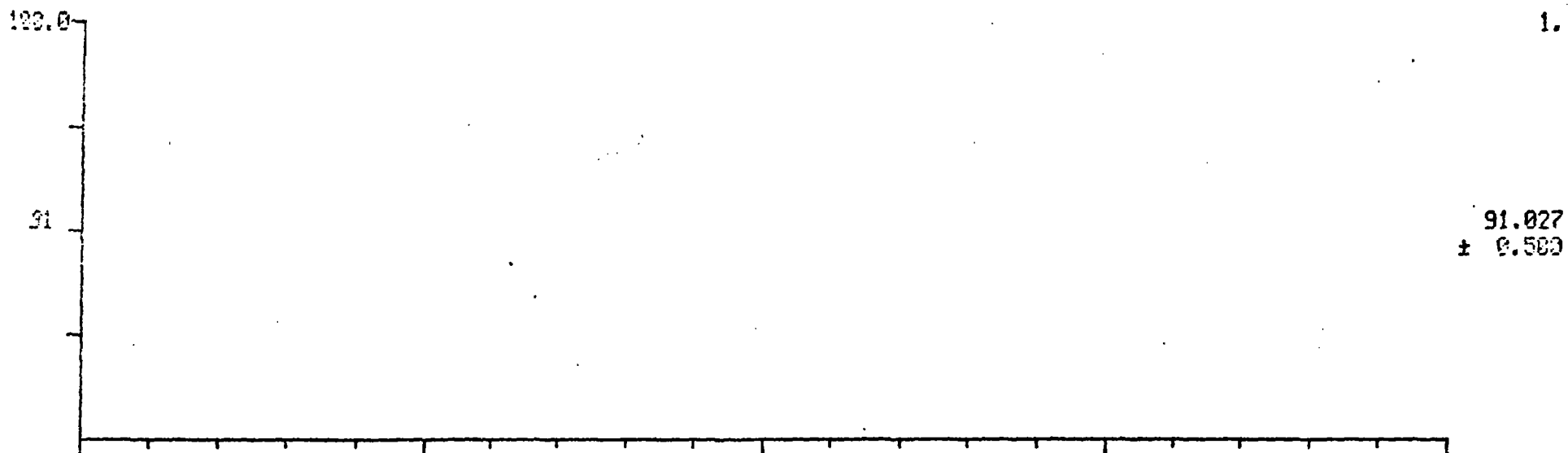
SCANS 539 TO 1000



MASS CHROMATOGRAMS
03/04/85 6:41:33
SAMPLE: BLANK HB1 CASE 3931 CH

DATA: V2715

SCANS 900 TO 1100

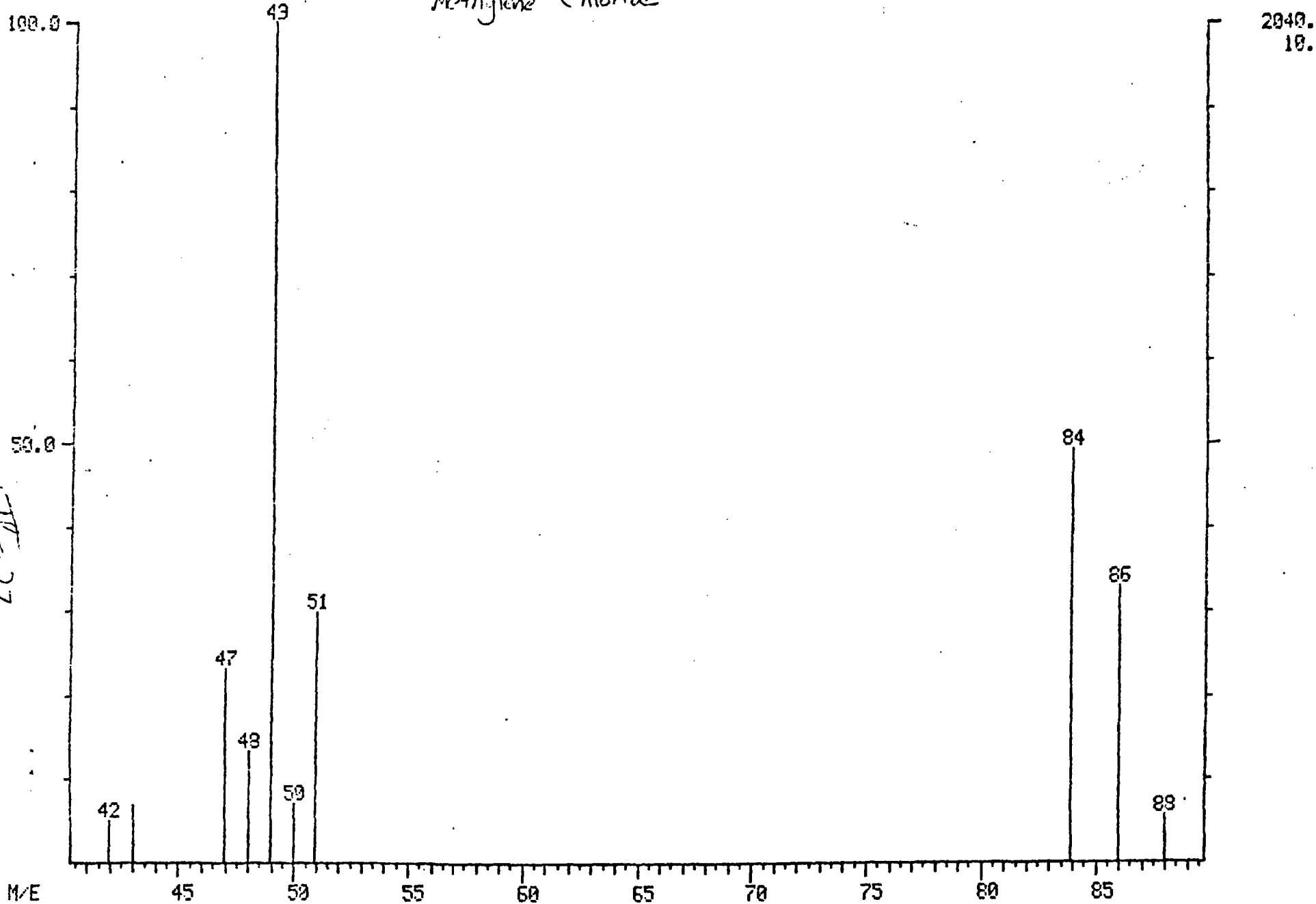


MASS SPECTRUM
00/04/85 6:41:00 + 8:03
SAMPLE: BLANK 181 CASE 3931 CM
ENHANCED (S 15B 2N)

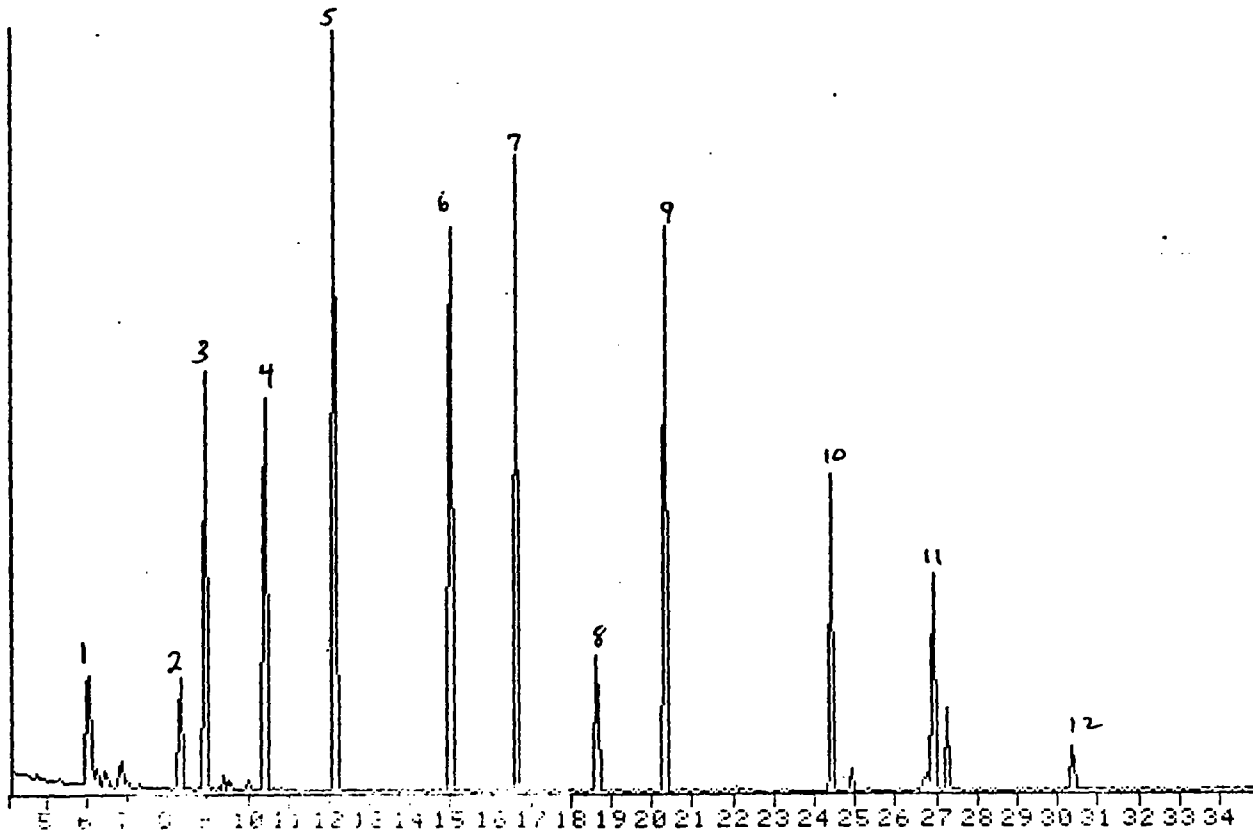
DATA: U2715 #253

BASE M/E: 49
RIC: 5576.

Methylene Chloride



TIA



1. 2-FLUOROPHENOL (SURR)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBIENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

QC-956 MB1 CASE 3969
 3/25/85/EM

FILE NUMBER 14756
 BTL#15 Q14756 D14756

LINE	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
0	150.0	1.000	80.0000	.0000		D4-1,4-DICHLORO BENZENE(20)	
9	74.0	.046	.0000	.0000		N-NITROSODIMETHYLAMINE	
4	93.0	.913	.0000	.0000		ANILINE	
5	93.0	.709	.0000	.0000		BIS(2-CHLOROETHYL)ETHER	
3	146.0	.591	.0000	.0000		1,3-DICHLORO BENZENE	
5	146.0	.594	.0000	.0000		1,4-DICHLORO BENZENE	
5	108.0	.395	.0000	.0000		BENZYL ALCOHOL	
5	146.0	.594	.0000	.0000		1,2-DICHLORO BENZENE	
4	45.0	.527	.0000	.0000		BIS(2-CHLOROISOPROPYL)ETHER	
2	70.0	.490	.0000	.0000		N-NITROSODINPROPYLAMINE	
5	117.0	.291	.0000	.0000		HEXACHLOROETHANE	
5	77.0	.716	.0000	.0000		NITROBENZENE	

1	136.0	1.000	80.0000	.0000		D8-NAPHTHALENE(20)	
1	138.0	.089	.0000	.0000		ISOPHORONE	
0	95.0	.106	.0000	.0000		BIS(2-CHLOROETHOXY)METHANE	
1	180.0	.154	.0000	.0000		1,2,4-TRICHLORO BENZENE	
0	129.0	.077	.0000	.0000		NAPHTHALENE	
1	127.0	.267	.0000	.0000		4-CHLOROANILINE	
1	225.0	.070	.0000	.0000		HEXACHLORO BUTADIENE	
0	115.0	.151	.0000	.0000		2-METHYLNAPHTHALENE	

10	68.0	1.000	80.0000	.0000		D8-NAPHTHALENE(20)	
11	138.0	1.519	.0000	.0000		ISOPHORONE	
11	95.0	1.813	.0000	.0000		BIS(2-CHLOROETHOXY)METHANE	
12	180.0	2.642	.0000	.0000		1,2,4-TRICHLORO BENZENE	
12	129.0	1.325	.0000	.0000		NAPHTHALENE	
12	127.0	4.577	.0000	.0000		4-CHLOROANILINE	
14	225.0	1.190	.0000	.0000		HEXACHLORO BUTADIENE	
14	115.0	2.578	.0000	.0000		2-METHYLNAPHTHALENE	

QC-956 MB1 CASE 3969
3/25/85/EM

FILE NUMBER 14756

BTL#15 Q14756

D14756

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
16.7	164.0	1.000	80.0000				D10-ACENAPHTHENE
14.6	237.0	.205	.0000				HEXACHLOROCYCLOPENTADIENE
15.2	164.0	.281	.0000				2-CHLORONAPHTHALENE
15.6	138.0	.397	.0000				2-NITROANILINE
16.2	164.0	.083	.0000				DIMETHYL PHTHALATE
16.3	151.0	.272	.0000				ACENAPHTHYLENE
16.4	165.0	.228	.0000				2,6-DINITROTOLUENE
16.7	138.0	.293	.0000				3-NITROANILINE
16.8	152.0	.437	.0000				ACENAPHTHENE
17.2	139.0	.490	.0000				DIBENZOFURAN
17.3	165.0	.320	.0000				2,4-DINITROTOLUENE
18.0	177.0	.171	.0000				DIETHYL PHTHALATE
18.0	165.0	.820	.0000				FLUORENE
18.1	204.0	.358	.0000				4-CHLOROPHENYLPHENYL ETHER
18.2	138.0	.292	.0000				4-NITROANILINE
17.2	169.0	.144	.0000				DIPHENYLAMINE
20.3	188.0	1.000	80.0000				D10-PHENANTHRENE(20)
19.3	248.0	.112	.0000				4-BROMOPHENYLPHENYL ETHER
19.6	284.0	.112	.0000				HEXACHLOROBENZENE
20.4	179.0	.060	.0000				PHENANTHRENE/ANTHRACENE
22.1	150.0	.100	.0000				DIBUTYL PHTHALATE
23.4	202.0	.665	.0000				FLUORANTHENE
26.9	240.0	1.000	80.0000				D12-CHRYSENE
23.9	202.0	2.264	.0000				PYRENE
25.8	149.0	1.136	.0000				BUTYL BENZYL PHTHALATE
26.9	228.0	.427	.0000				BENZO(A)ANTHRACENE/CHRYSENE
27.3	149.0	1.128	24.0288				BIS(2-ETHYLHEXYL)PHTHALATE
30.4	264.0	1.000	80.0000				D12-PERYLENE
28.7	149.0	5.974	.0000				DIOCTYL PHTHALATE
29.5	252.0	1.307	.0000				BENZO(B)/(K)FLUORANTHENE
30.2	252.0	.744	.0000				BENZO(A)PYRENE
33.7	278.0	.485	.0000				INDENO(1,23-CD)PYRENE
33.8	278.0	.346	.0000				DIBENZO(AH)ANTHRACENE
34.6	270.0	.360	.0000				BENZO(GHI)PERYLENE

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FILE NUMBER 14756

QC-956 MB1 CASE 3969

3/25/85/EM

BTL#15 Q14756

D14756

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
26.9	240.0	1.000	80.0000				D12-CHRYSENE
23.8	184.0	.712	.0000				BENZIDINE
27.0	252.0	.146	.0000				DICHLOROBENZIDINE

IV-31

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
9.0	150.0	1.000	80.0000				D4-1,4-DICHLOROBENZENE(20)
8.4	94.0	.841	.0000				PHENOL
8.6	128.0	.611	.0000				2-CHLOROPHENOL
9.9	108.0	.671	.0000				2-METHYLPHENOL
10.2	108.0	.699	.0000				4-METHYLPHENOL

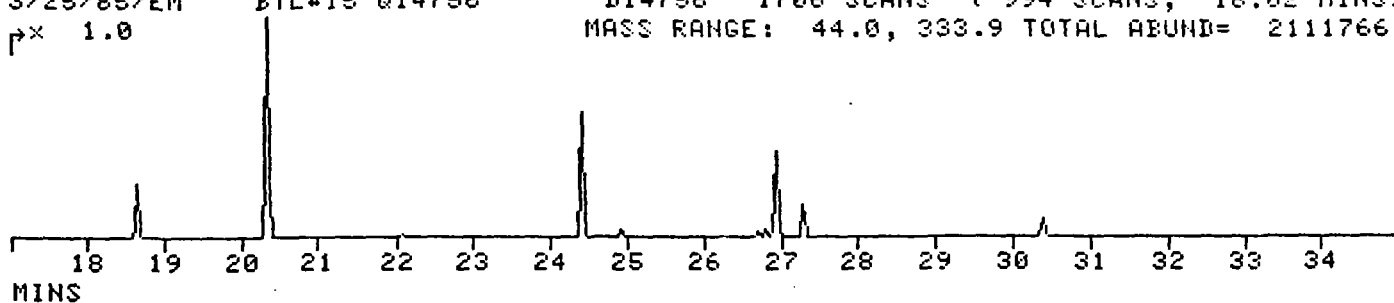
12.1	68.0	1.000	80.0000				D8-NAPHTHALENE(20)
11.3	139.0	1.759	.0000				2-NITROPHENOL
11.5	122.0	2.581	.0000				2,4-DIMETHYLPHENOL
12.1	122.0	1.200	.0000				BENZOIC ACID
11.9	162.0	2.051	.0000				2,4-DICHLOROPHENOL
13.8	142.0	2.196	.0000				4-CHLORO-M-CRESOL

16.7	164.0	1.000	80.0000				D10-ACENAPHTHENE(20)
14.8	196.0	.237	.0000				2,4,5-TRICHLOROPHENOL
14.8	196.0	.237	.0000				2,4,6-TRICHLOROPHENOL
17.0	184.0	.111	.0000				2,4-DINITROPHENOL
18.3	198.0	.126	.0000				4,6-DINITRO-O-CRESOL
17.2	65.0	.231	.0000				4-NITROPHENOL

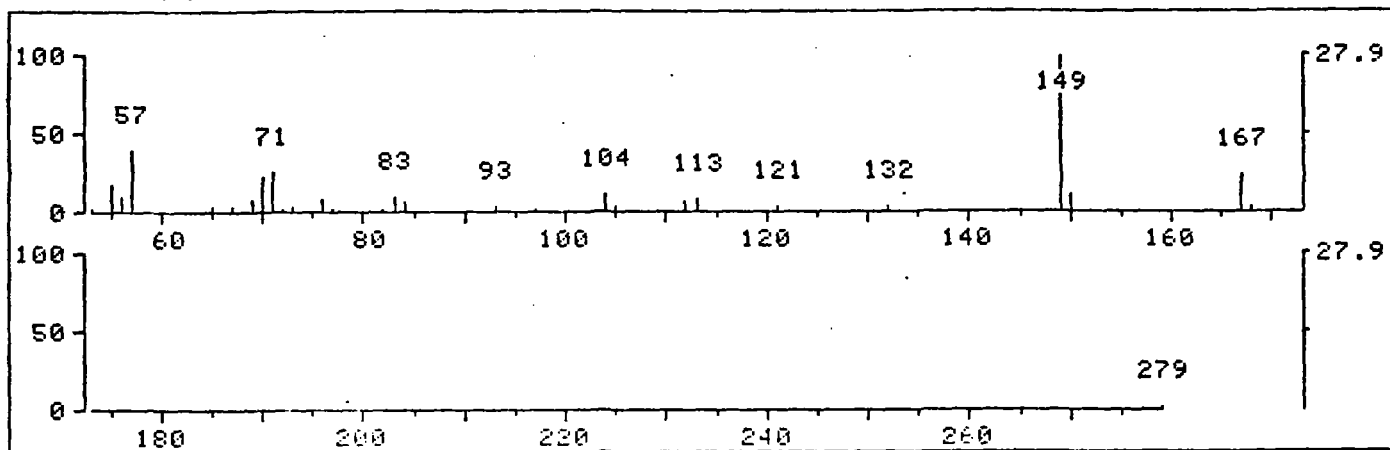
20.3	188.0	1.000	80.0000				D10-PHENANTHRENE
20.1	266.0	.072	.0000				PENTACHLOROPHENOL

QC-956 MB1 CASE 3969
3/25/85/EM BTL#15 Q14756
p x 1.0

D14756 1706 SCANS (994 SCANS, 18.02 MINS)
MASS RANGE: 44.0, 333.9 TOTAL ABUND= 2111766.

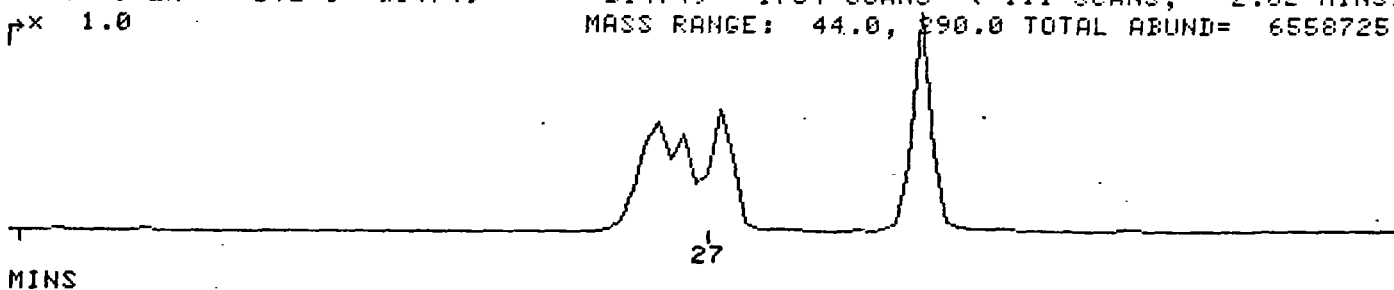


*1280 RET. TIME: 27.28 TOT ABUND= 6671. BASE PK/ABUND: 149.1/ 2415.

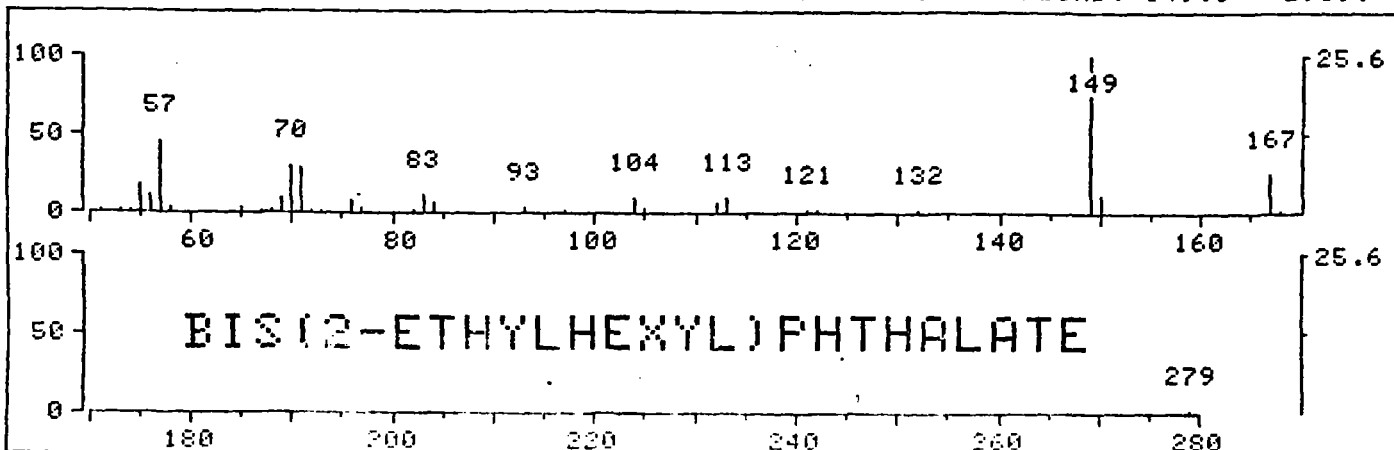


808N
3/25/85/EM BTL#8 Q14749
p x 1.0

D14749 1704 SCANS (111 SCANS, 2.02 MINS)
MASS RANGE: 44.0, 390.0 TOTAL ABUND= 6558725.



*1278 RET. TIME: 27.28 TOT ABUND= 11554. BASE PK/ABUND: 149.0/ 2959.



HEX TABLE ENTRIES: FRN 14756

Entry	Time	Mass	Area	%
1	27.0	239.7	14878.	100.0
2	24.9	TI	3902.	26.2

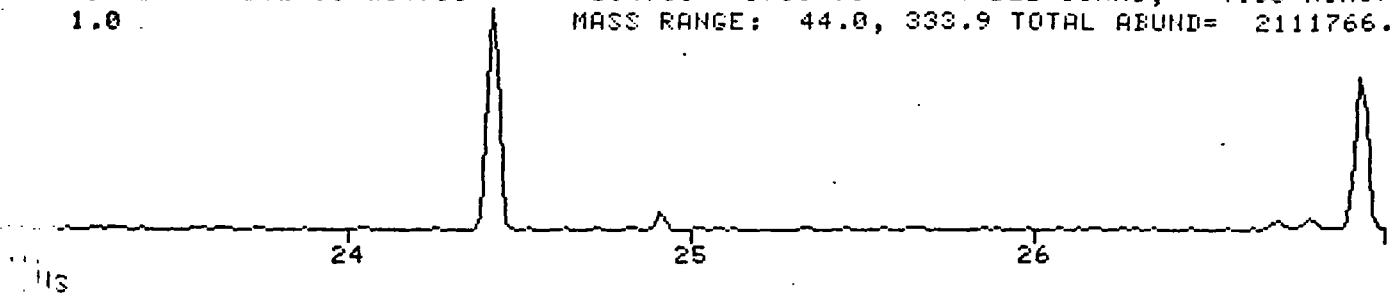
CALCULATE % ON ENTRY #:

V-34

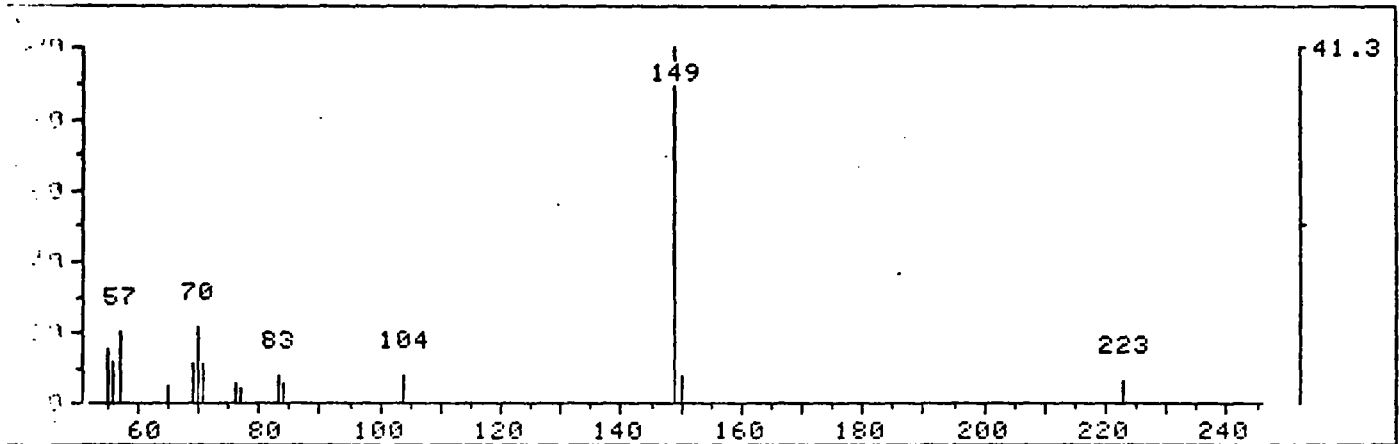
206

956 MB1 CASE 3969
5/85/EM BTL#15 Q14756
1.0

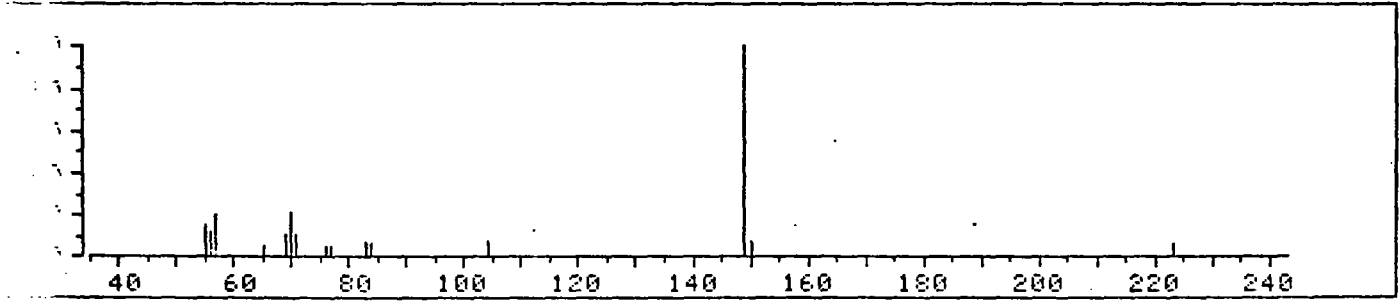
FRN 14756, SCAN 7
D14756 1706 SCANS (222 SCANS, 4.03 MINS)
MASS RANGE: 44.0, 333.9 TOTAL ABUND= 2111766.



1149 RET. TIME: 24.90 TOT ABUND= 2692. BASE PK/ABUND: 149.1/ 1113.

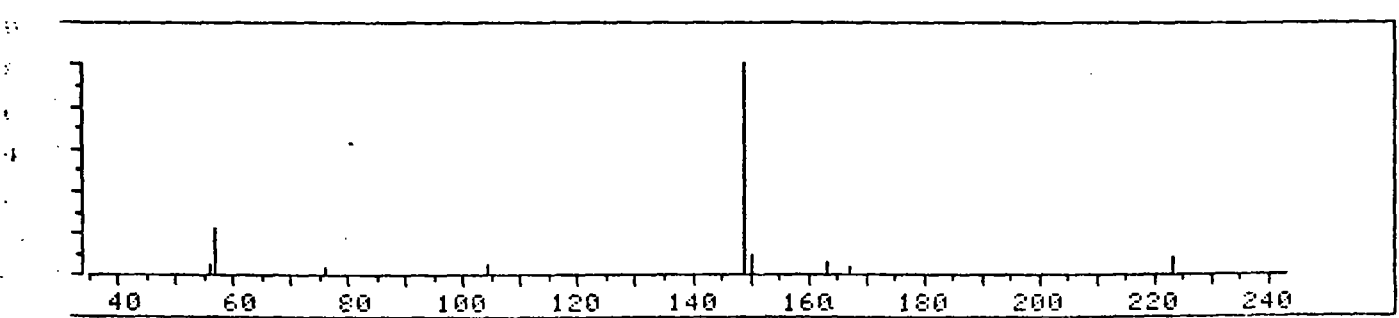


HITS: REFERENCE FRN 14756 SCAN 1149



.7%

6 LFRN 3011 SPECT 803 MW= 278 C16H22O4
59 1,2-Benzenedicarboxylic acid, bis(2-methylpropyl) ester (9CI)



V-35

24

Date 3/30/75

Analyst BL

Contract 3-705-050

Pesticide Analysis - Qualitative/Quantitative

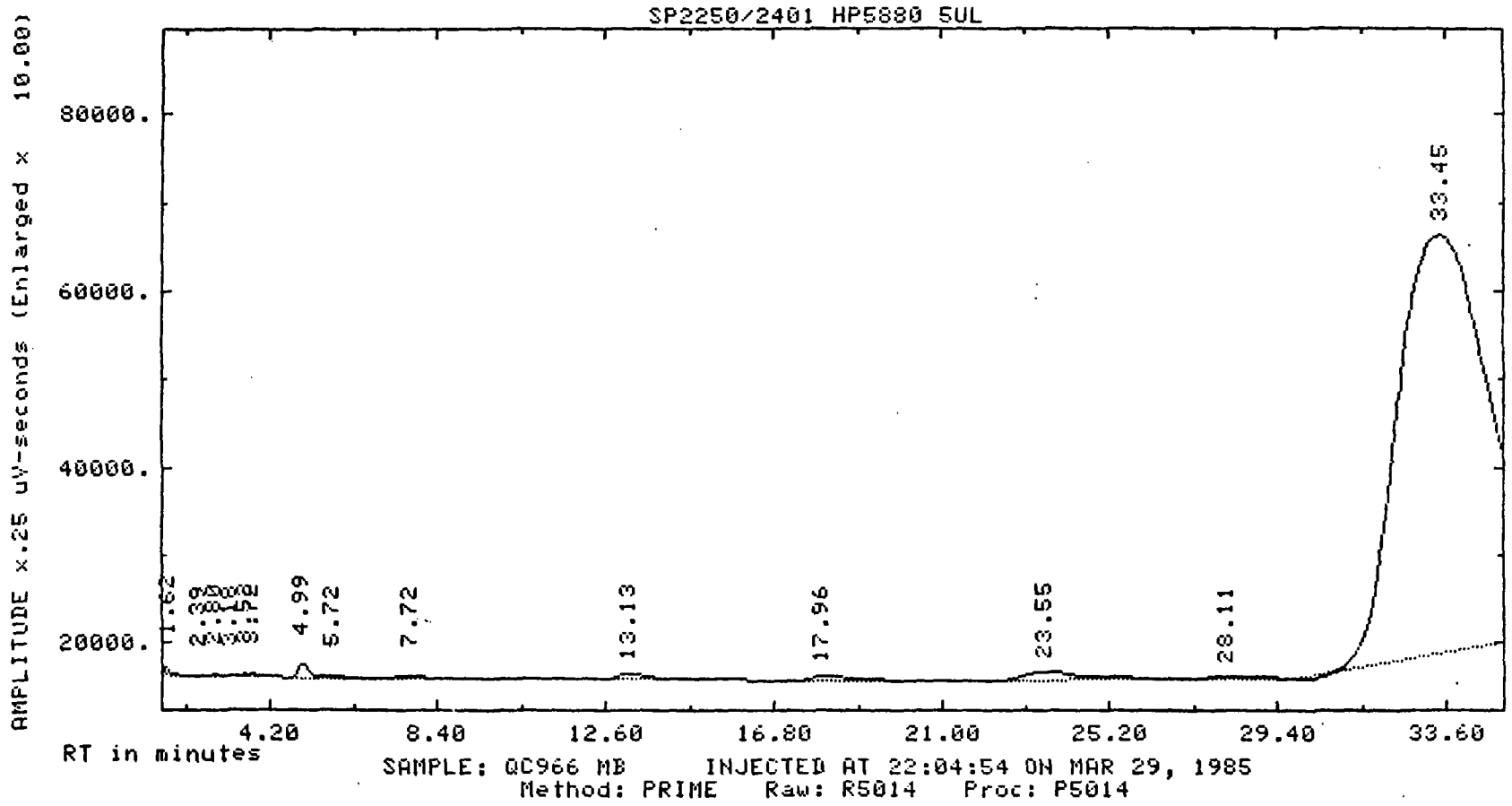
Case # 3969

Sample I.D. QC-966 / Method Blank 1
Run # 14

HP 58 80
Bottle # 14

RT (min.)	Peak ID	Area x 10 ³	ng/ml	Dilution Factor (mls extracted)	(ug/l) ppb	total ug
3.18	Heptachlor	0.24	<5.0	x100 x10 ÷ 1000 =	<0.05	
3.52	δ-BHC	1.26	<5.0	x10 ÷ 1000 =	<0.05	
	DBC obscured by peak at 33 min					

Case 3969



V-37

APR 2, 1985 11:35:23

REPORT: 110. CHANNEL: 2 # PEAKS: -18 SP2250/2401 HP5880 5UL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.02	0.00	1.00000	347 BV	.007	
2	1.31	0.00	1.00000	144 VB	.003	
3	1.62	0.00	1.00000	3618 BB	.076	
4	2.39	0.00	1.00000	203 BB	.004	
5	2.80	0.00	1.00000	1605 BV	.035	
6	3.18	0.00	1.00000	241 VV	.005	
7	3.52	0.00	1.00000	1258 VV	.026	
8	3.72	0.00	1.00000	1664 VB	.035	
9	4.99	0.00	1.00000	14528 BV	.308	
10	5.72	0.00	1.00000	1810 VB	.038	
11	7.72	0.00	1.00000	7321 BB	.154	
12	13.12	0.00	1.00000	16297 BB	.343	
13	17.98	0.00	1.00000	18845 BB	.396	
14	23.55	0.00	1.00000	42044 BV	.884	
15	28.11	0.00	1.00000	10597 VV	.349	
16	33.45	0.00	1.00000	4183503 VV	88.001	
17	37.18	0.00	1.00000	303871 VV	7.654	
18	38.99	0.00	1.00000	79887 VF	1.600	

TOTAL AREA = 4754062 TOTAL AREA % = 100.000

SAMPLE: Q080E MB INJECTED AT 22:04:54 ON MAR 29, 1985

ZERO METHOD: PRIME SEQ: SMO3 SUBSEQ/SAMP: 1 / 14 BTL: 14

SL-WDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	NO

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

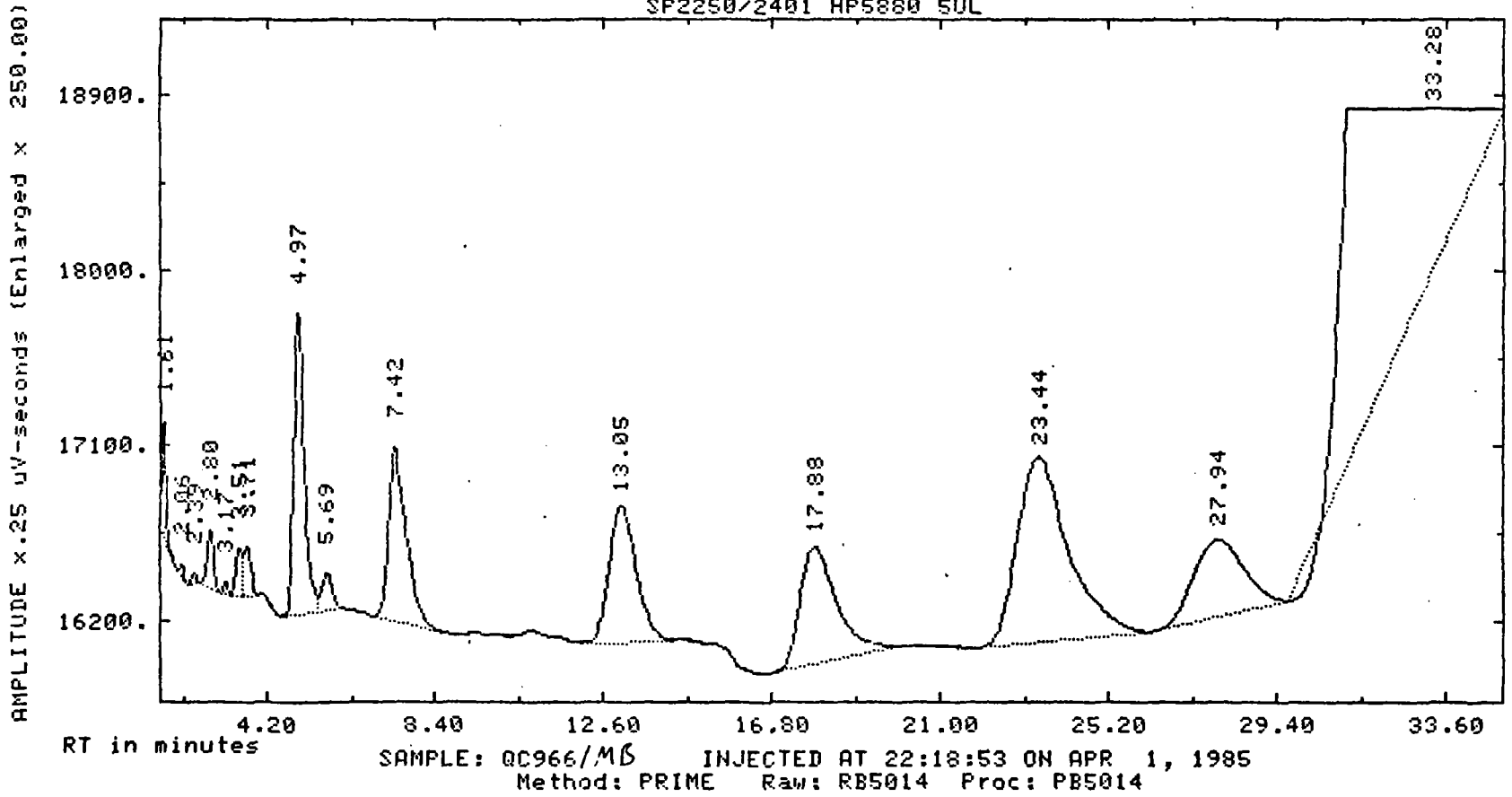
ENDED NOT ON BASELINE

RAW DATA FILE: R5014 PARAM FILES= METHOD: SEQ:

DONE
READY

Case 3969

SP2250/2401 HP5880 5UL



68-1

*LI,P,PB5014

PROCESSED DATA FILE: PB5014 ON CRN 21

Case 3969

APR 2, 1985 9:13:17

REPORT: 157 CHANNEL: 2 # PEAKS: -18 SP2250/2401 HP5890 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.01	0.00	1.00000	276 BV	.007	
2	1.31	0.00	1.00000	235 VB	.006	
3	1.61	0.00	1.00000	2063 BB	.051	
4	2.06	0.00	1.00000	160 BB	.004	
5	2.39	0.00	1.00000	180 BB	.004	
6	2.80	0.00	1.00030	1553 BV	.039	
7	3.17	0.00	1.00000	240 UV	.006	
8	3.51	0.00	1.00000	1250 UV	.031	
9	3.71	0.00	1.00000	1577 VB	.039	
10	4.97	0.00	1.00000	14158 BV	.349	
11	5.69	0.00	1.00000	1760 VB	.043	
12	7.42	0.00	1.00000	13370 BB	.330	
13	13.05	0.00	1.00000	16151 BB	.398	
14	17.80	0.00	1.00000	16848 BB	.465	
15	23.44	0.00	1.00000	41699 BV	1.029	
16	27.94	0.00	1.00000	16789 UV	.414	
17	33.28	0.00	1.00000	3800389 UV	95.366	
18	39.29	0.00	1.00000	57522 VF	1.419	

TOTAL AREA = 4054240 TOTAL AREA % = 100.000

SAMPLE: Q0966/MB INJECTED AT 22:19:53 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 14 BTL: 14

SL-WIDTH	MV/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: RB5014 PARAM FILES= METHOD: SEQ:

DONE
READY

II-40

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Organics Analysis Data Sheet
(Page 1)

Laboratory Name: GCA
 Lab Sample ID No. Blank
 Sample Matrix Water
 Data Release Authorized By _____

Case No. 3931
 QC Report No. 3931
 Contract No. 68-01-6267
 Date Sample Received 3/5/85

Volatiles Compounds

Concentration Low Medium (Circle One)
 Date Extracted/Prepared 3/5/85
 Date Analyzed 3/5/85
 Conc/Dil Factor NA pH NA
 Percent Moisture NA
 Percent Moisture (Decanted) NA

CAS Number	Compound	ug/l or ug/kg (Circle One)
74-87-3	Chloroethane	10u
74-83-9	Bromoethane	10u
75-01-4	Vinyl Chloride	10u
75-00-3	Chloroethane	10u
75-48-2	Methylene Chloride	2.3
67-64-1	Acetone	10u
75-15-0	Carbon Disulfide	5u
75-35-4	1,1-Dichloroethane	5u
75-34-3	1,1-Dichloroethane	5u
100-10-5	Trans-1,2-Dichloroethane	5u
67-66-3	Chloroform	5u
107-06-2	1,2-Dichloroethane	5u
75-53-5	2-Butanone	10u
71-65-6	1,1,1-Trichloroethane	5u
75-23-5	Carbon Tetrachloride	5u
108-01-4	Vinyl Acetate	10u
71-27-4	Bromochloroethane	5u

CAS Number	Compound	ug/l or ug/kg (Circle One)
75-34-5	1,1,2,2-Tetrachloroethane	5u
78-87-5	1,2-Dichloropropane	5u
100-10-6	Trans-1,3-Dichloropropane	5u
79-01-6	Trichloroethane	5u
124-36-1	Dibromochloromethane	5u
79-00-5	1,1,2-Trichloroethane	5u
71-41-7	Benzene	5u
100-10-5	cis-1,3-Dichloropropane	5u
110-25-8	2-Chloroethylvinylether	10u
75-25-7	Bromoform	5u
591-78-6	2-Hexanone	10u
108-10-1	4-Methyl-2-Pentanone	10u
127-18-4	Tetrachloroethane	5u
108-28-3	Toluene	5u
108-90-7	Chlorobenzene	5u
100-41-4	Fluorobenzene	5u
100-42-5	Styrene	5u
	Total Volatiles	

Value: Indicates the value of the compound detected. If the value is less than the detection limit, the value is reported as ND.

U: Indicates the value of the compound detected. If the value is less than the detection limit, the value is reported as ND. If the value is greater than the detection limit, the value is reported as U. If the value is greater than the detection limit, the value is reported as U. If the value is greater than the detection limit, the value is reported as U.

D: Indicates the value of the compound detected. If the value is less than the detection limit, the value is reported as ND. If the value is greater than the detection limit, the value is reported as D. If the value is greater than the detection limit, the value is reported as D.

Other: Other compounds or parameters may be reported by the analyst. If so, they must be fully identified and their values reported to the data user.

7-11 II-41

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	<i>No Compounds Found</i>	<i>Y01A</i>		
2.				
3.				
4.				
5.				
6.				
7.				
8.				
9.				
10.				
11.				
12.				
13.				
14.				
15.				
16.				
17.				
18.				
19.				
20.				
21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				

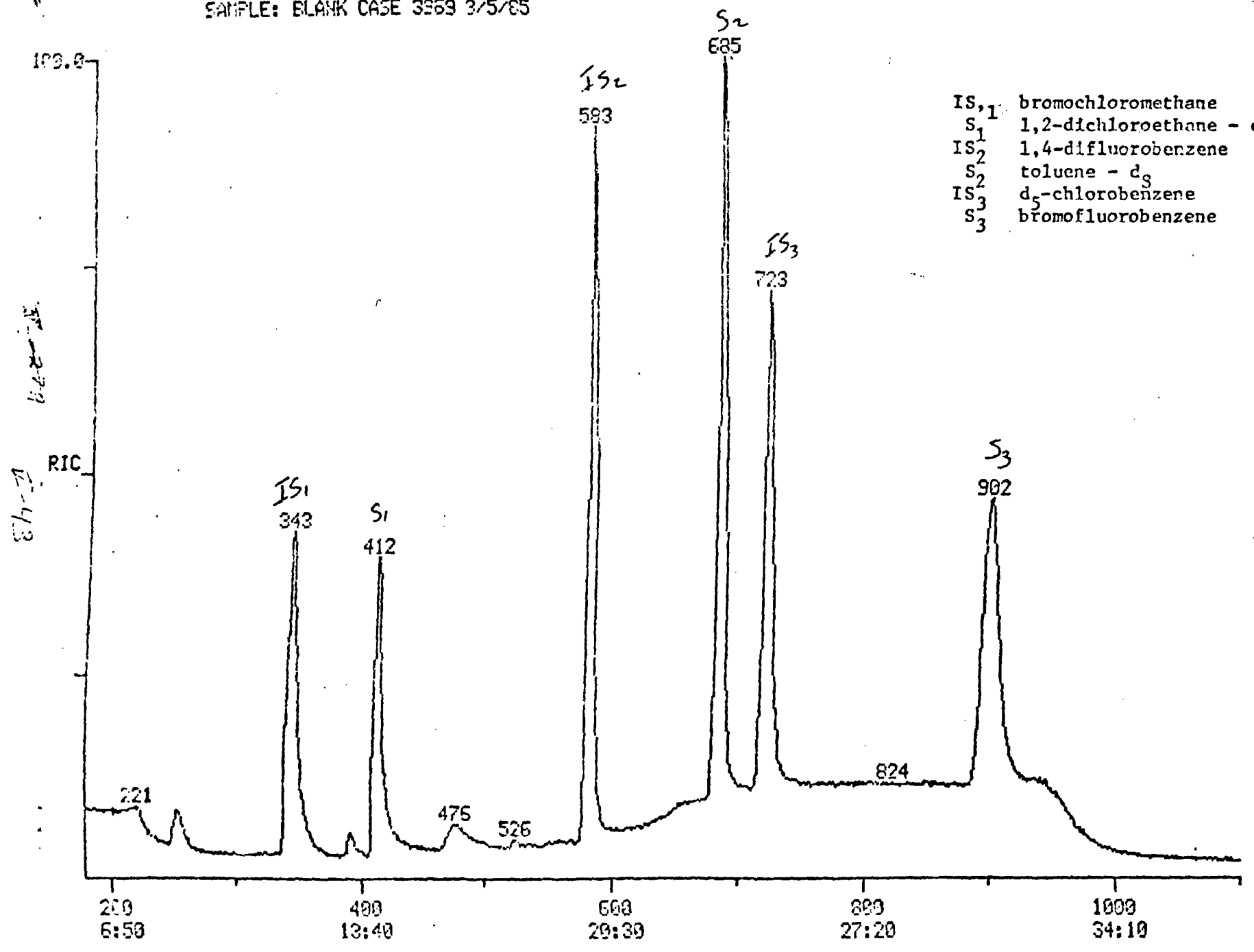
*7-78
T.C.H.*

RIC
03/05/85 9:51:00
SAMPLE: BLANK CASE 3363 3/5/85

DATA: U2735

SCANS 180 TO 1100

112769.

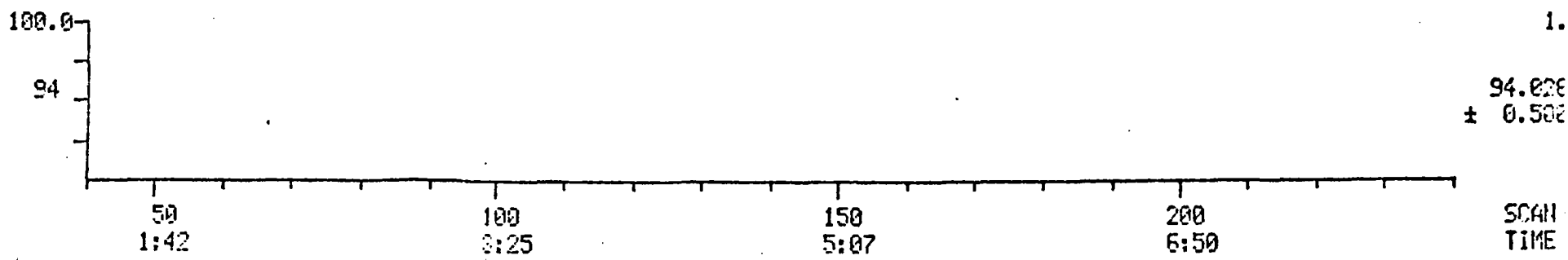
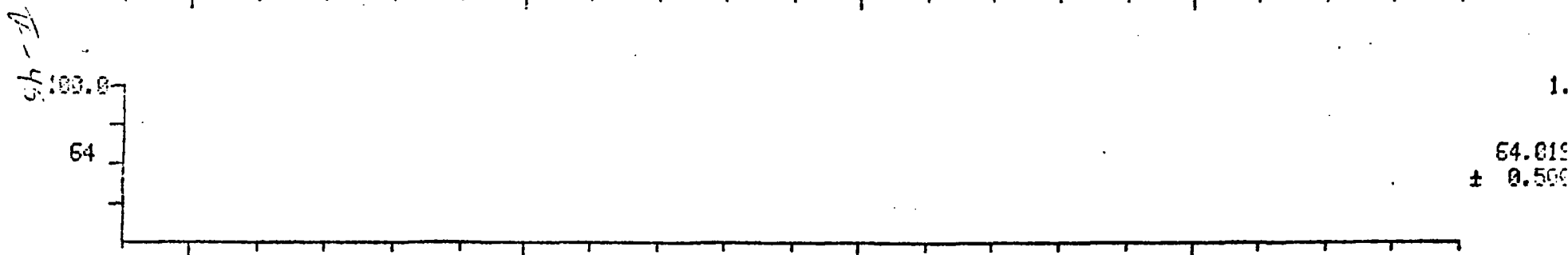
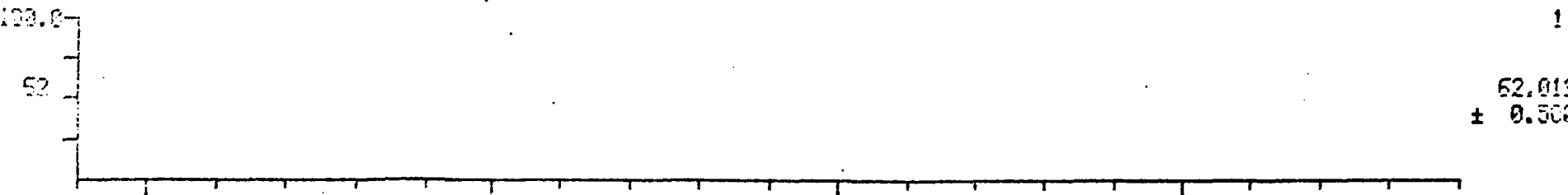
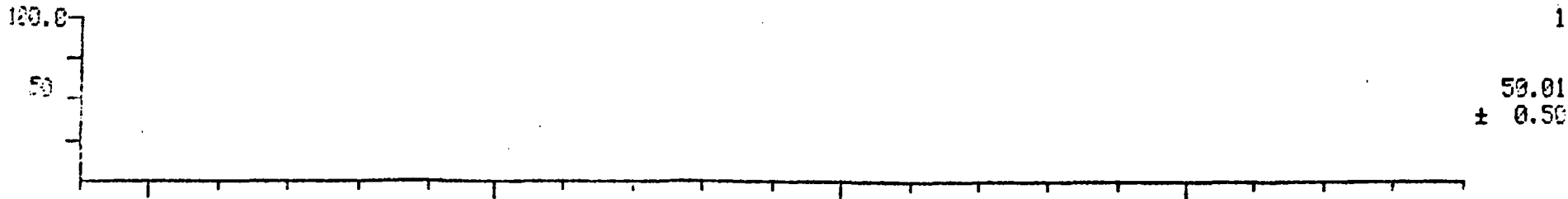


Case: 3969

NO	NAME
1	BROMOCHLOROMETHANE (I.S. #1)
2	CHLOROMETHANE
3	BROMOMETHANE
4	VINYL CHLORIDE
5	CHLOROETHANE
6	METHYLENE CHLORIDE
7	CARBON DISULFIDE
8	ACETONE
9	1,1-DICHLOROETHENE
10	1,1-DICHLOROETHANE
11	TRANS-1,2-DICHLOROETHENE
12	CHLOROFORM
13	D4-1,2-DICHLOROETHANE (SURR #1)
14	1,2-DICHLOROETHANE
15	1,1,1-TRICHLOROETHANE
16	CARBON TETRACHLORIDE
17	VINYL ACETATE
18	2-BUTANONE
19	1,4-difluorobenzene (I.S. #2)
20	BROMODICHLOROMETHANE
21	1,2-DICHLOROPROPANE
22	TRANS-1,3-DICHLOROPROPENE
23	TRICHLOROETHENE
24	DIBROMOCHLOROMETHANE
25	1,1,2-TRICHLOROETHANE
26	CIS-1,2-DICHLOROPROPENE
27	BENZENE
28	2-CHLOROETHYL VINYL ETHER
29	BROMOFORM
30	D8-TOLUENE
31	D5-CHLOROBENZENE (I.S. #3)
32	TETRACHLOROETHENE
33	1,1,2,2-TETRACHLOROETHANE
34	2-METHYL-2-PENTANONE
35	2-HEXANONE
36	TOLUENE
37	CHLOROBENZENE
38	ETHYL BENZENE
39	BFB (SURR #3)
40	STRYENE
41	TOTAL XYLENES

44

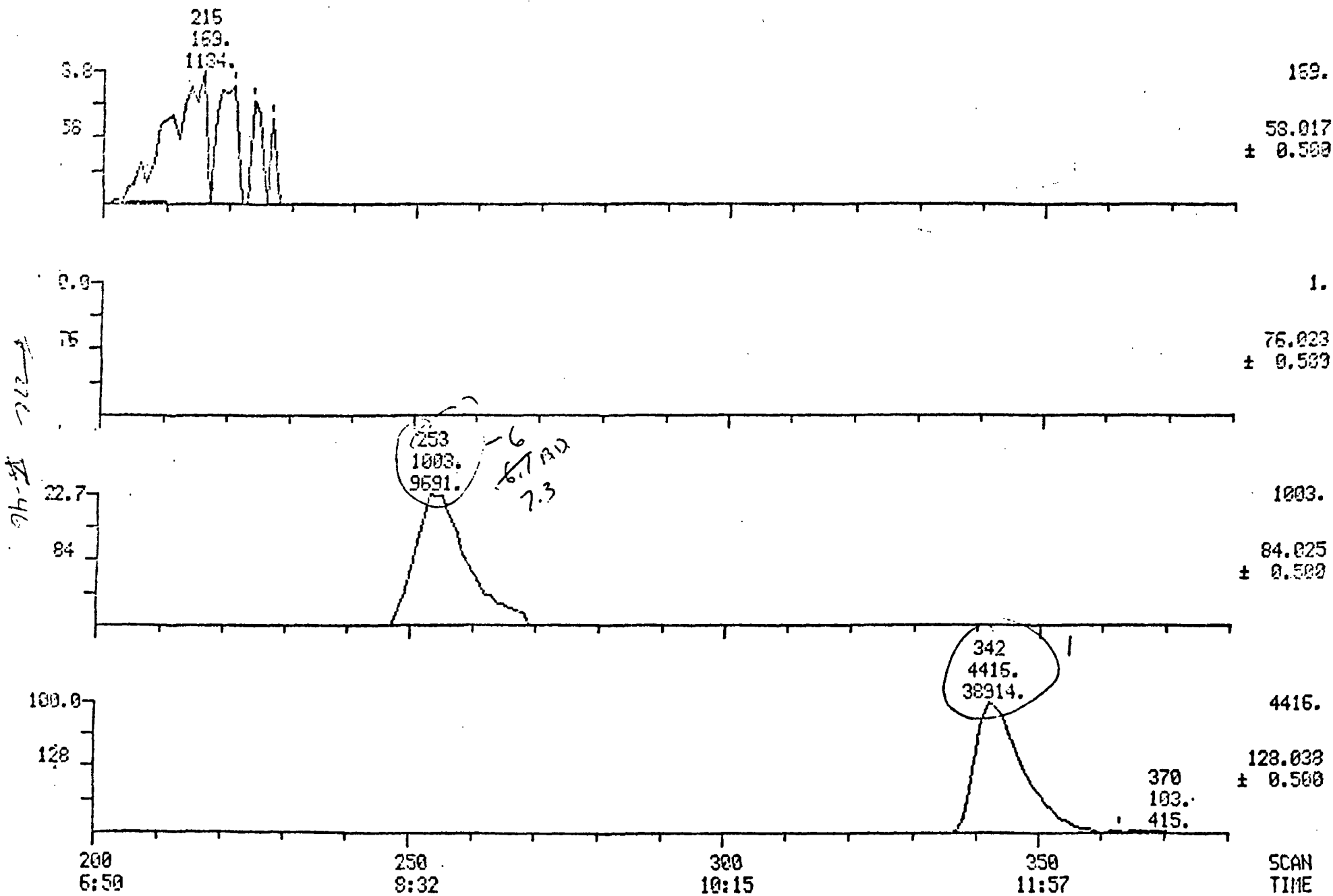
03/05/85 9:51:00
SAMPLE: BLANK CASE 3959 3/5/85



MASS CHROMATOGRAMS
03/05/85 9:51:00
SAMPLE: BLANK CASE 3969 3/5/85

DATA: U2735

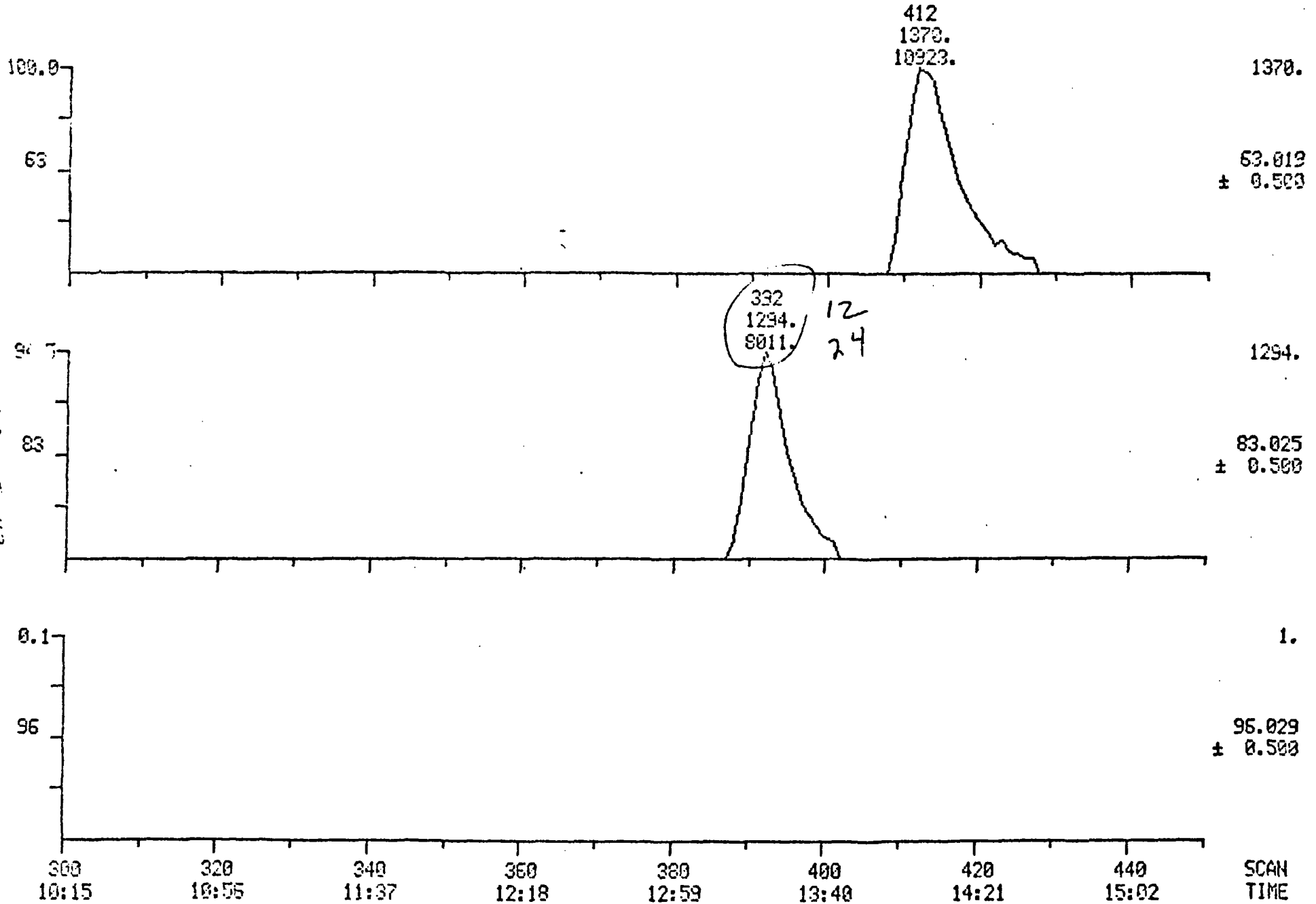
SCANS 200 TO 380



MASS CHROMATOGRAMS
03/05/85 9:51:30
SAMPLE: BLANK CASE 3963 3/5/85

DATA: V2736

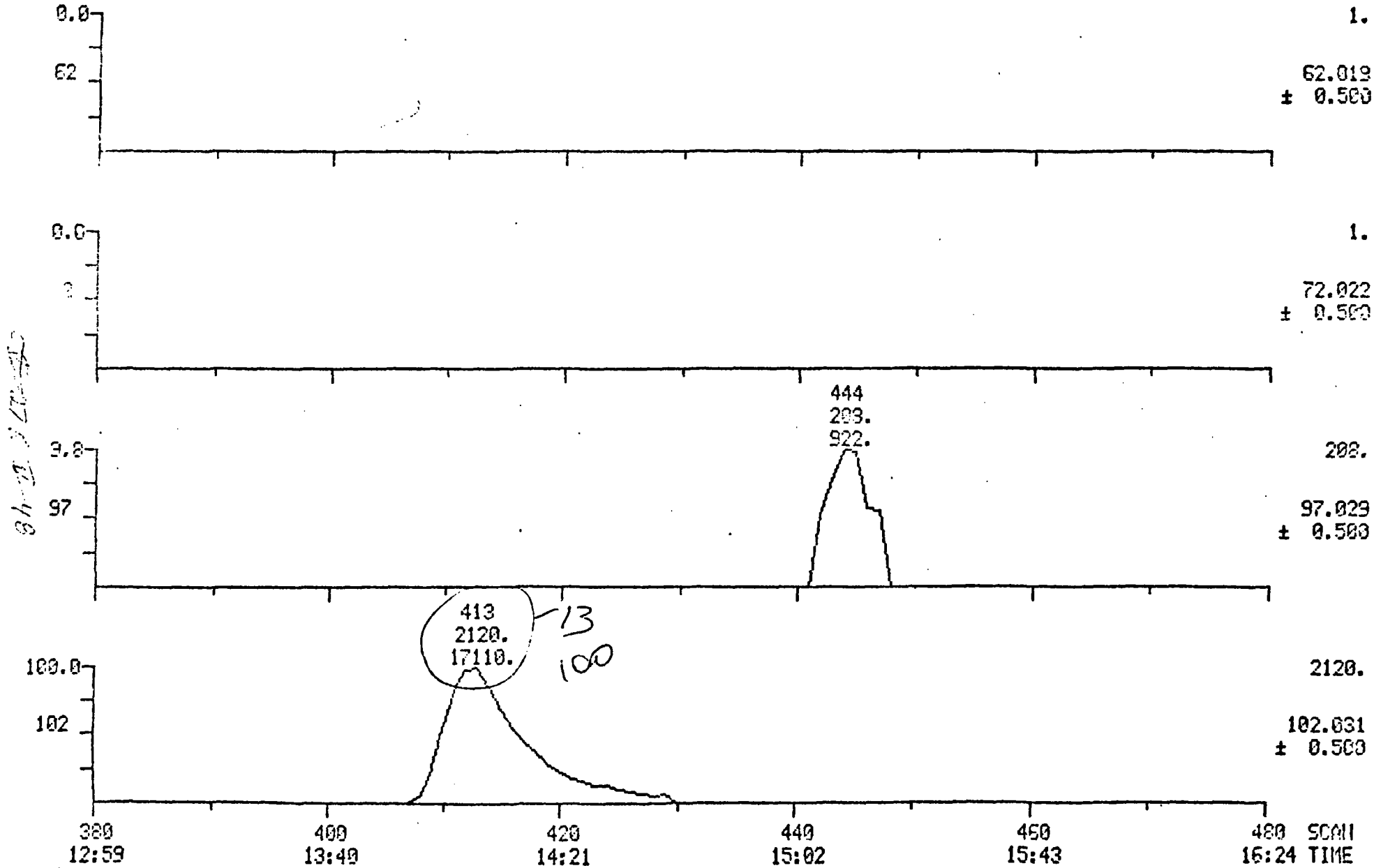
SCANS 300 TO 450



MASS CHROMATOGRAMS
03/05/85 9:51:00
SAMPLE: BLANK CASE 3959 3/5/85

DATA: U2736

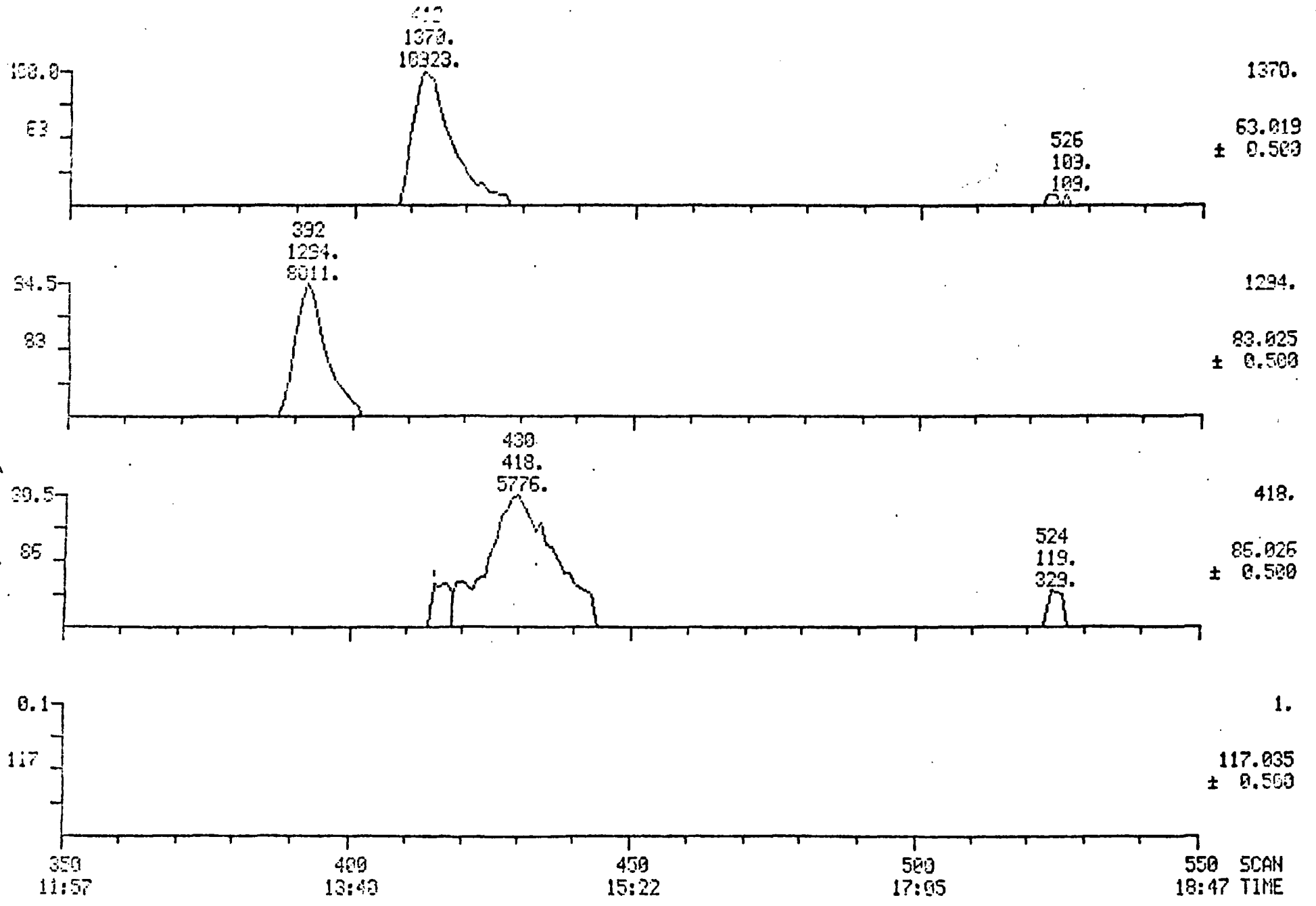
SCANS 380 TO 480



MASS CHROMATOGRAMS
03/25/85 9:51:00
SAMPLE: BLANK CASE 3969 3/5/85

DATA: U2736

SCANS 350 TO 550



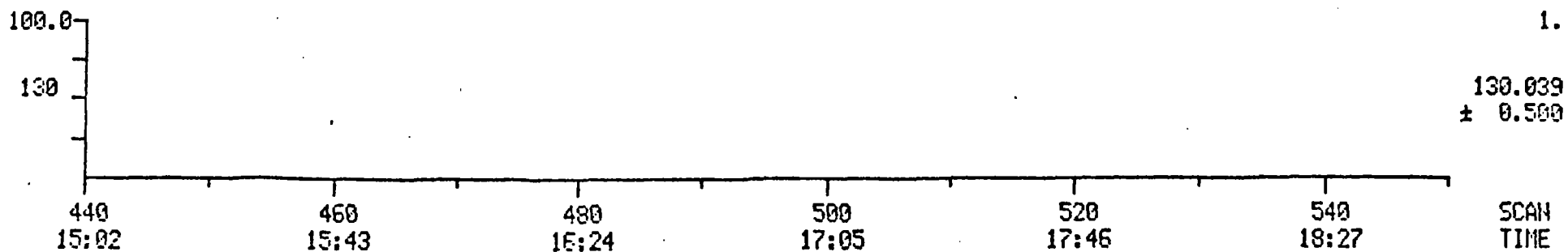
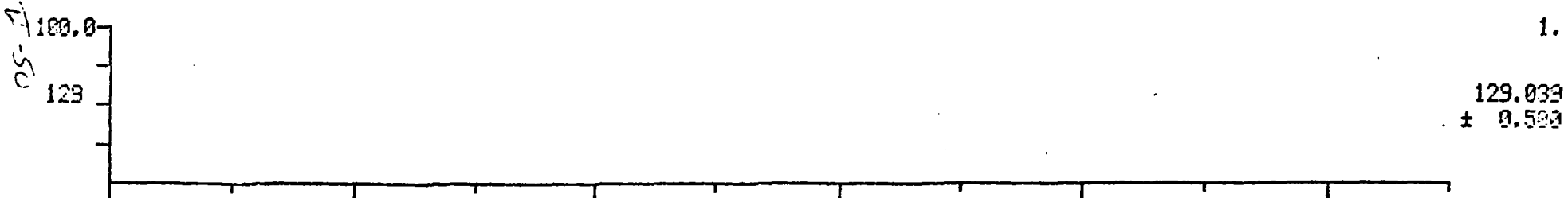
blank injection

MS

MASS CHROMATOGRAMS
03/25/85 9:51:00
SAMPLE: BLANK CASE 3959 3/5/85

DATA: U2736

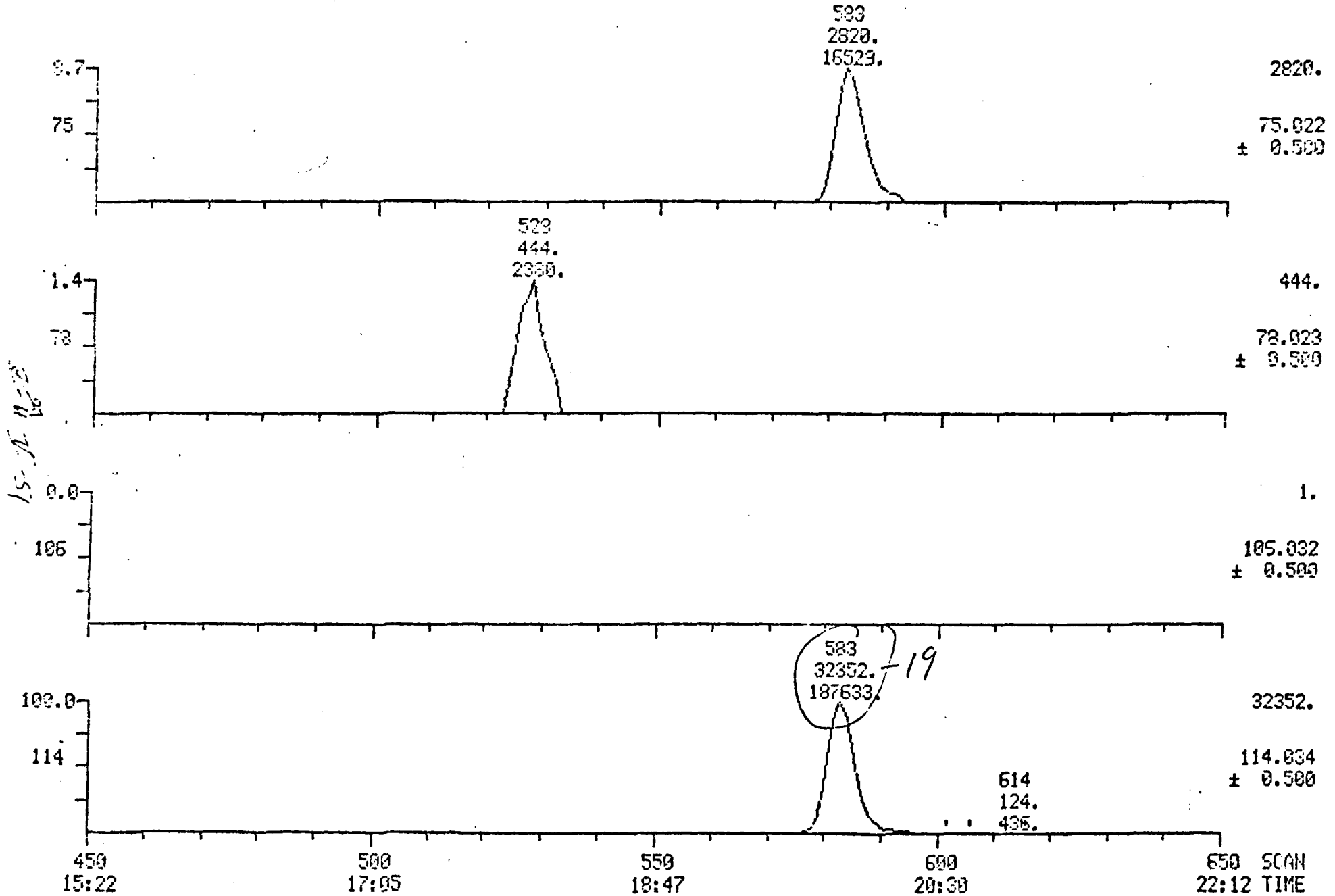
SCANS 440 TO 550



MASS CHROMATOGRAMS
03/05/85 9:51:00
SAMPLE: BLANK CASE 3953 3/5/85

DATA: U2736

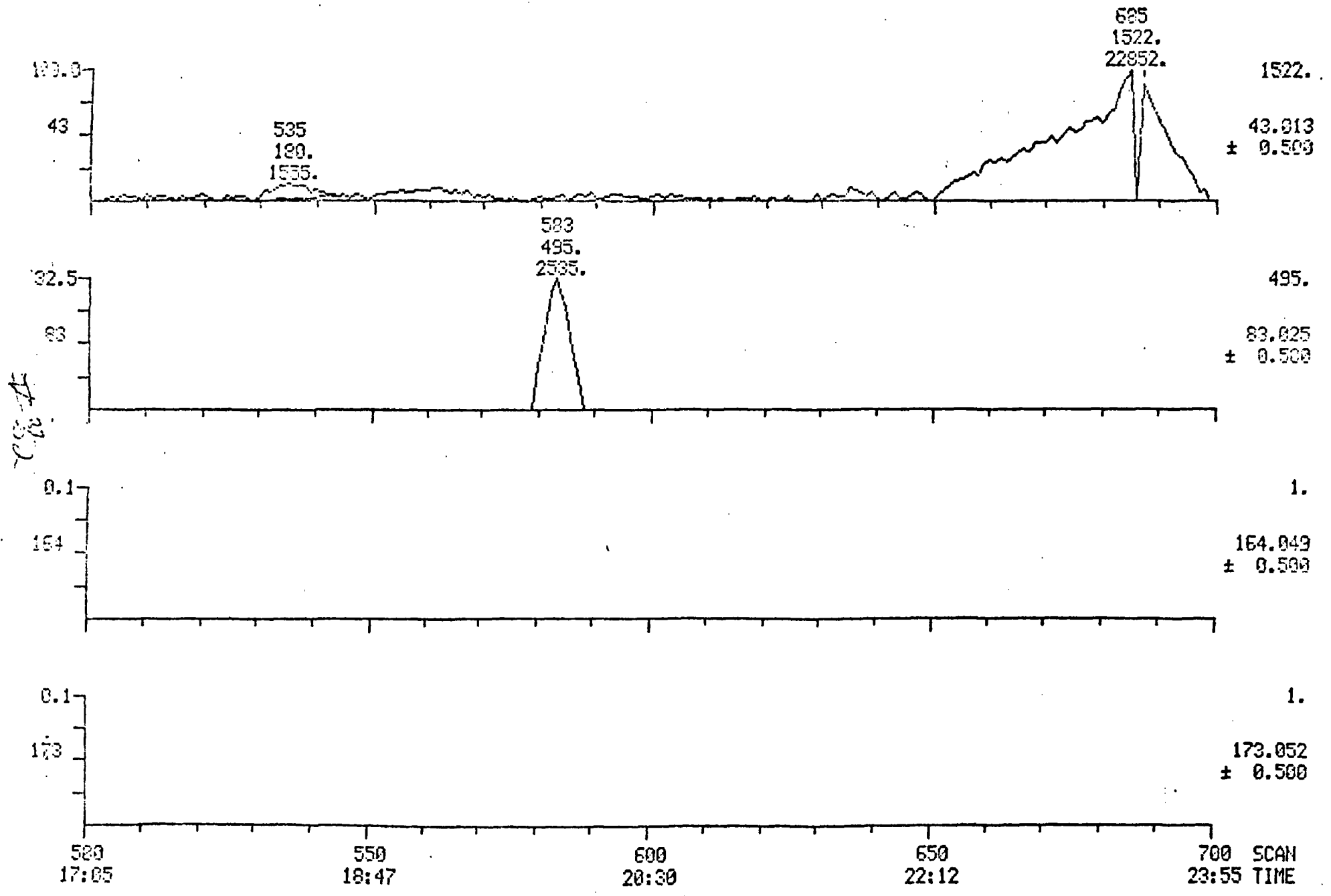
SCANS 450 TO 650



MASS CHROMATOGRAMS
03/03/85 9:01:00
SAMPLE: BLANK CASE 3969 3/5/85

DATA: U2735

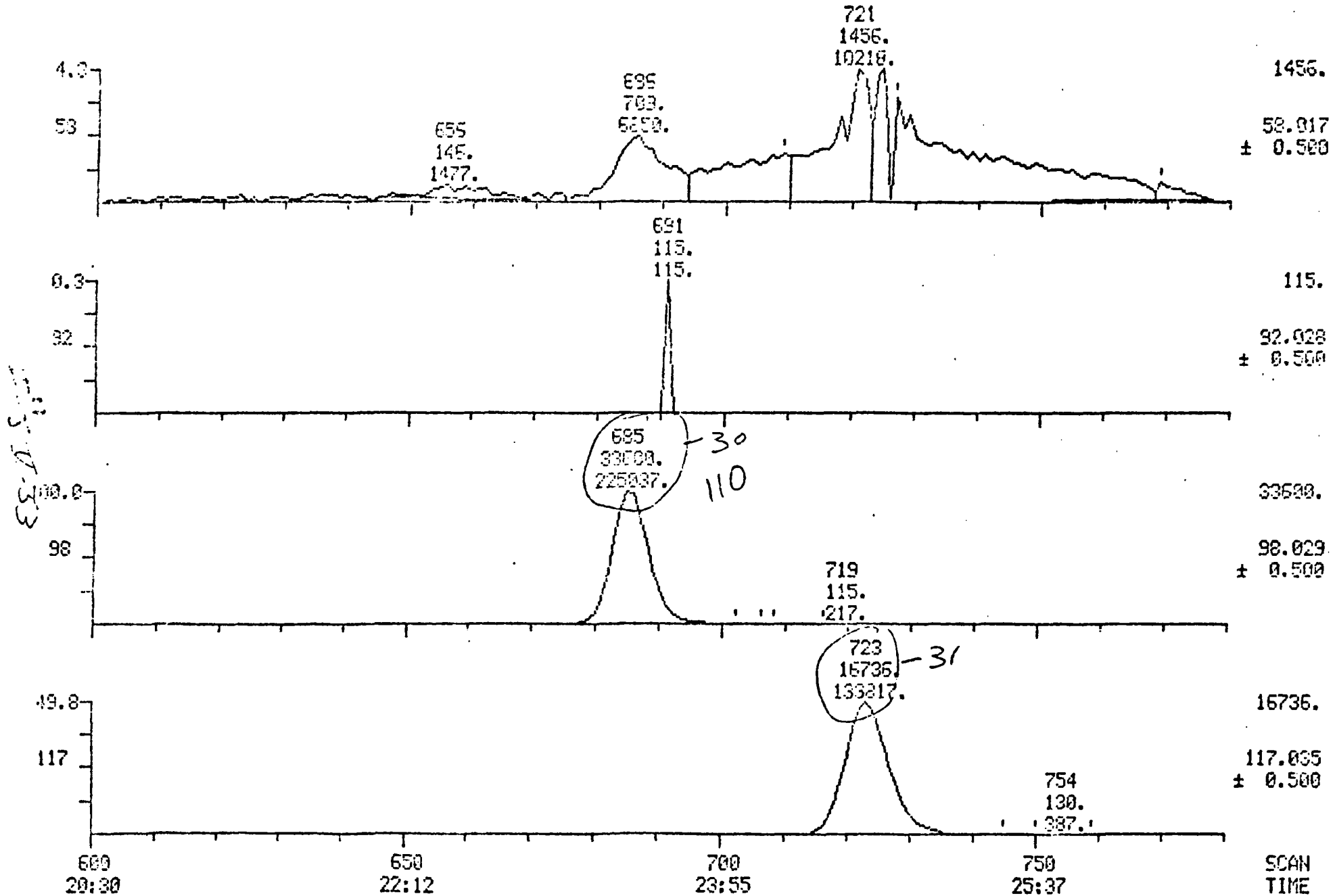
SCANS 500 TO 700



MASS CHROMATOGRAMS
03/05/85 9:51:00
SAMPLE: BLANK CASE 3369 3/5/85

DATA: U2736

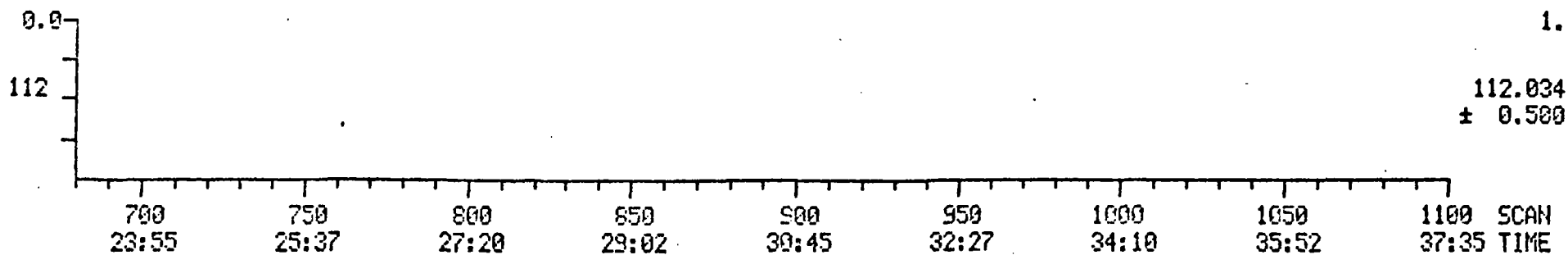
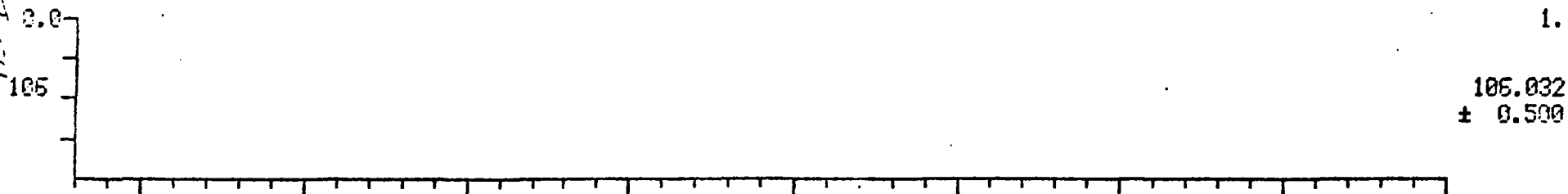
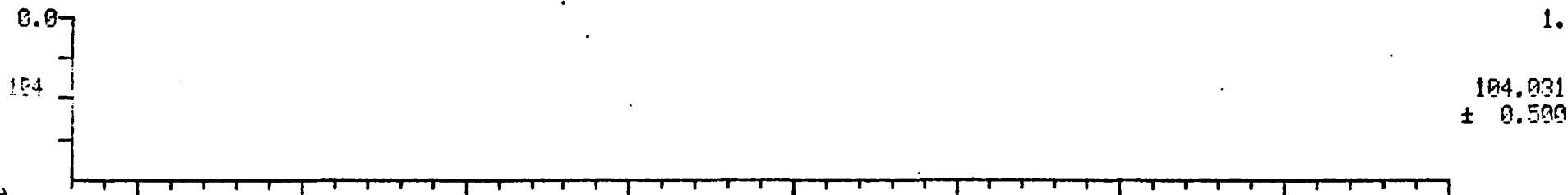
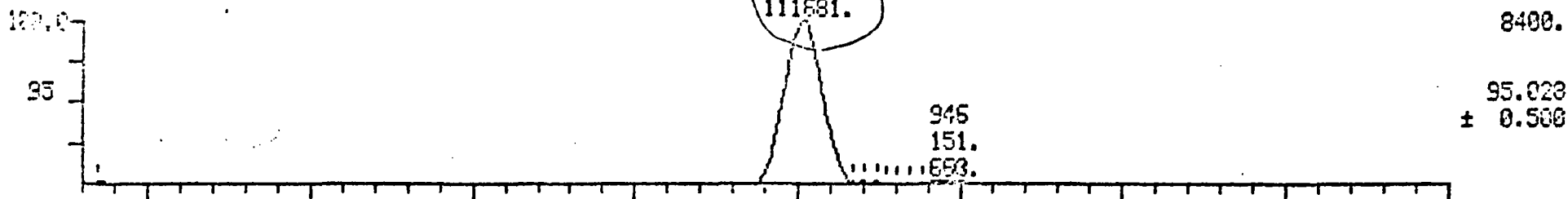
SCANS 603 TO 789



MASS CHROMATOGRAMS
03/05/95 9:51:00
SAMPLE: BLANK CASE 3969 3/5/85

DATA: U2736

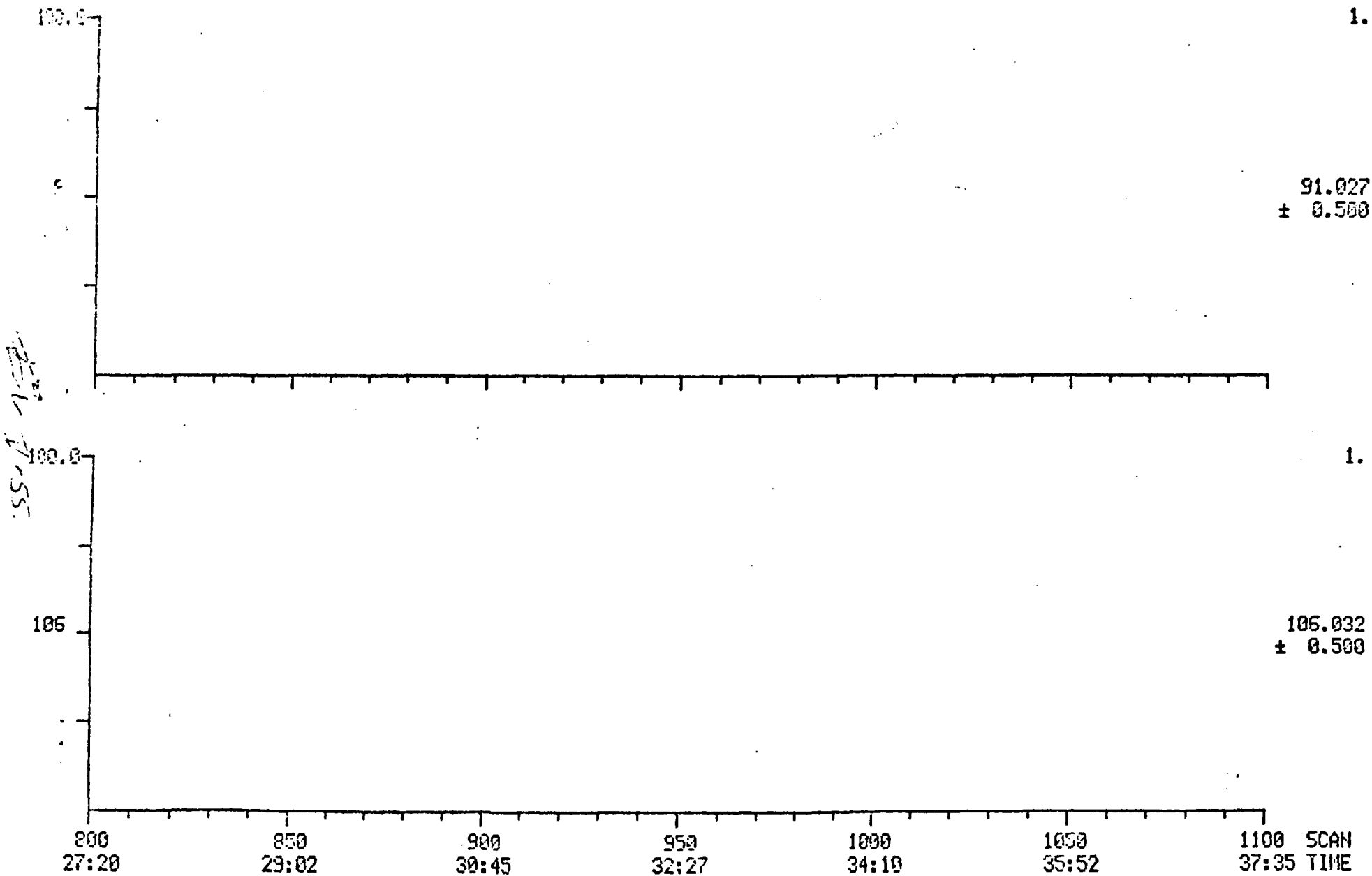
SCANS 680 TO 1100



MASS CHROMATOGRAMS
03/05/85 9:51:00
SAMPLE: BLANK CASE 3569 3/5/85

DATA: U2736

SCANS 800 TO 1100



Laboratory Name: GCA
Lab Sample ID No: blak
Sample Matrix: water
Data Release Authorized By: _____

Case No: 3969
QC Report No: 3969
Contract No: 68-01-6767
Date Sample Received: 3/6/85

Volatile Compounds

Concentration Low Medium (Circle One)
Date Extracted/Prepared: 3/6/85
Date Analyzed: 3/6/85
Conc/Dil Factor: 1 pH: NA
Percent Moisture: NA
Percent Moisture (Decanted): NA

CAS Number	Compound	ug/L or ug/Kg (Circle One)
74-87-3	Chloromethane	10u
74-83-9	Bromomethane	10u
75-01-4	Vinyl Chloride	10u
75-00-3	Chloroethane	10u
75-00-2	Ethyl Chloride	2.5
67-64-1	Ethylene	10u
75-15-0	Carbon Dioxide	5u
75-35-4	1,1-Dichloroethane	5u
75-34-3	1,2-Dichloroethane	5u
116-07-5	Trans-1,2-Dichloroethane	5u
07-55-3	Chloroform	5u
107-06-2	1,2-Dichlorobenzene	5u
76-53-9	Trichloroethane	10u
71-55-6	1,1,1-Trichloroethane	5u
35-28-5	Carbon Tetrachloride	5u
106-06-4	o-Cresol	10u
78-27-4	Bromochloroethane	5u

CAS Number	Compound	ug/L or ug/Kg (Circle One)
79-34-5	1,1,2,2-Tetrachloroethane	5u
78-97-5	1,2-Dichloropropane	5u
10061-02-6	trans-1,3-Dichloropropene	5u
72-01-7	Trichloroethane	5u
124-49-1	Dibromochloromethane	5u
79-00-5	1,1,2-Trichloroethane	5u
71-43-2	Benzene	5u
10061-01-5	cis-1,3-Dichloropropene	5u
110-75-8	2-Chloroethylchloride	10u
75-25-2	Bromobenzene	5u
501-75-0	2-Hexanone	10u
145-10-1	4-Methyl-2-Pentanone	10u
127-18-4	Tetrahydrofuran	5u
108-98-3	Toluene	5u
109-50-7	Chlorobenzene	5u
100-41-4	Hexibenzene	5u
110-42-5	Styrene	5u
	Total Xylenes	5u

Quality Assurance

For every sample analyzed, the following quality assurance measures were taken:
A. All analytical results were checked against the original data. However, the accuracy of the results cannot be guaranteed.

Value: If the result is a value greater than the applicable detection limit, then the value is reported.

U: Indicates concentration was a value greater than the detection limit. Report the concentration value and the sample ID. If the concentration is less than the detection limit, then the result is reported as "not detected". This is not the case if the detection limit is less than the detection limit. If the detection limit is less than the detection limit, then the result is reported as "not detected". The number in the parentheses indicates the detection limit for the sample.

J: Indicates that the sample was analyzed for a specific compound. The detection limit for this compound is indicated in parentheses. If the concentration is less than the detection limit, then the result is reported as "not detected". If the concentration is greater than the detection limit, then the result is reported as "detected". The number in the parentheses indicates the detection limit for the sample.

C: This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single compound pesticides (100 ug/L) in the final extract should be identified by GC/MS.

P: This flag indicates that the sample was analyzed for a specific compound. The detection limit for this compound is indicated in parentheses. If the concentration is less than the detection limit, then the result is reported as "not detected". If the concentration is greater than the detection limit, then the result is reported as "detected". The number in the parentheses indicates the detection limit for the sample.

Other: Other specific flags and detection limits may be required to properly define the results. If any of these flags are used, they must be fully defined in the final report. If any of these flags are used, they must be fully defined in the final report.

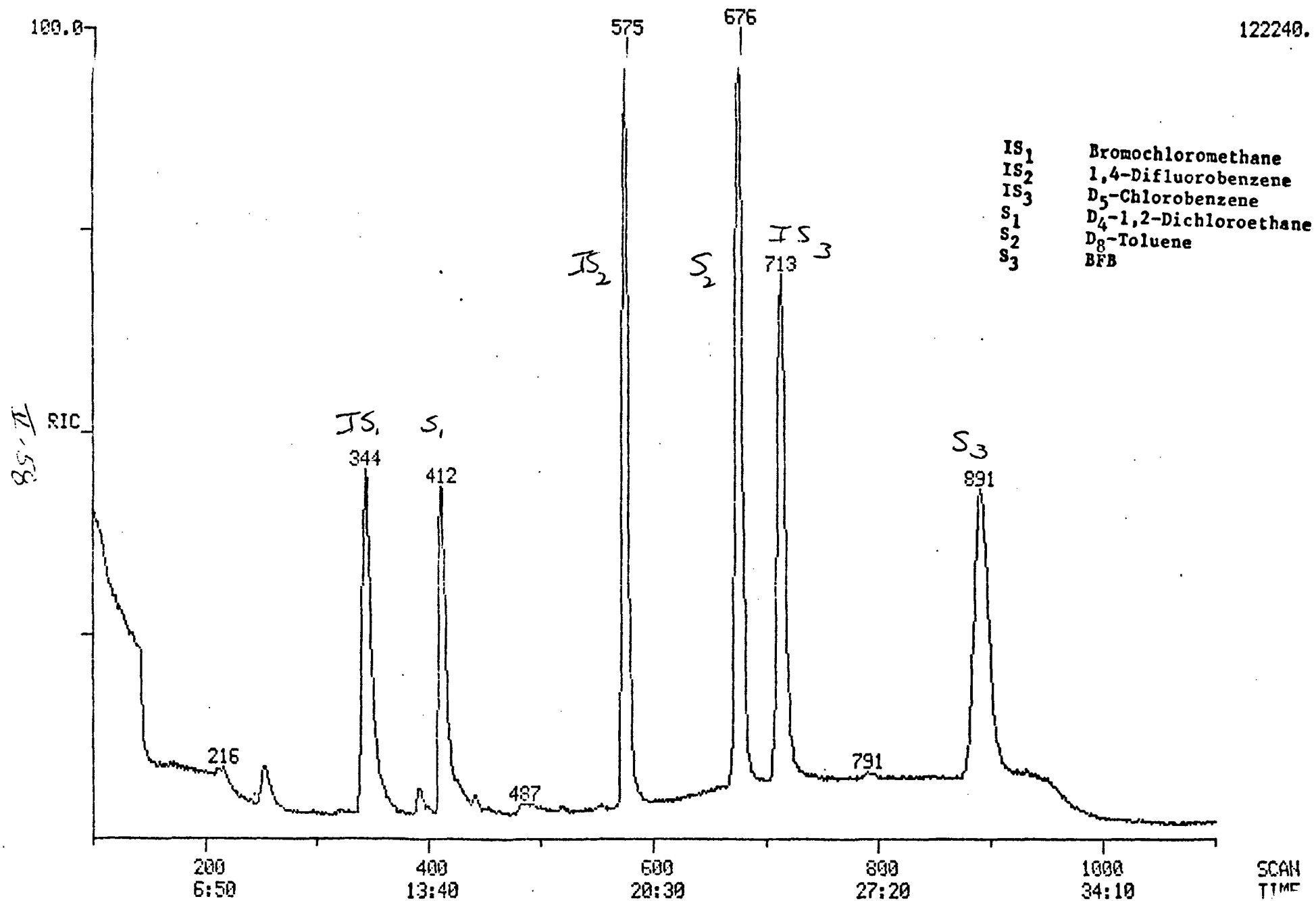
V-56

RIC
03/05/85 10:53:00
SAMPLE: BLANK 3/6/85 Case 3969

DATA: U2754

SCANS 100 TO 1100

122240.



06/85 10:53:00
IDENTIFICATION REPORT

FILE: M2

A: V2754.TI
06/85 10:53:00
FILE: BLANK 3/6/85
SUBMITTED BY:

MB3 Case 3969

ANALYST:

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

NAME
BROMOCHLOROMETHANE (I. S. #1)
CHLOROMETHANE
BROMOMETHANE
VINYL CHLORIDE
CHLOROETHANE
METHYLENE CHLORIDE
ACETONE
CARBON DISULFIDE
1, 1-DICHLOROETHENE
1, 1-DICHLOROETHANE
TRANS 1, 2-DICHLOROETHENE
CHLOROFORM
1, 2-DICHLOROETHANE
D4-1, 2-DICHLOROETHANE
1, 1, 1-TRICHLOROETHANE
1, 4-DIFLUOROBENZENE (I. S. #2)
CARBON TETRACHLORIDE
VINYL ACETATE
2-BUTANONE
BROMODICHLOROMETHANE
1, 2-DICHLOROPROPANE
TRANS 1, 3-DICHLOROPROPANE
TRICHLOROETHENE
DIBROMOCHLOROMETHANE
1, 1, 2-TRICHLOROETHANE
BENZENE
CIS-1, 3-DICHLOROPROPENE
BROMOFORM
D8-TOLUENE (SURR#2)
2-CHLOROETHYL VINYLETHER
D5-CHLOROBENZENE (I. S. #3)
4 METHYL 2 PENTANONE
2-HEXANONE
TETRACHLOROETHENE
1, 1, 2, 2, TETRACHLOROETHANE
TOLUENE
CHLOROBENZENE
ETHYLBENZENE
BFB (SURR#3)
STYRENE
TOTAL XYLENES

M/E	SCAN	TIME	REF	RRT	METH	AREA (HGHT)	AMOUNT	%TOT
128	344	11:45	1	1.000	A BB	43944.	50.000 UG/L	11.38
NOT FOUND								
NOT FOUND								

II-59

22

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
4		NOT FOUND							
5		NOT FOUND							
6		NOT FOUND							
7		NOT FOUND							
8		NOT FOUND							
9		NOT FOUND							
10		NOT FOUND							
11		NOT FOUND							
12		NOT FOUND							
13		NOT FOUND							
14	65	412	14:05	1	1.198	A BB	140000.	91.744 %REC	20.89
15		NOT FOUND							
16	114	575	19:39	16	1.000	A BB	222337.	50.000 UG/L	11.38
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		NOT FOUND							
27		NOT FOUND							
28		NOT FOUND							
29	98	675	23:04	16	1.174	A BB	254942.	96.660 %REC	22.00
30		NOT FOUND							
31	117	713	24:22	31	1.000	A BB	154346.	50.000 UG/L	11.38
32		NOT FOUND							
33		NOT FOUND							
34		NOT FOUND							
35		NOT FOUND							
36	92	681	23:16	31	0.955	A BB	1634.	0.471 UG/L	0.11
37		NOT FOUND							
38		NOT FOUND							
39	95	892	30:29	31	1.251	A BB	128970.	100.391 %REC	22.85
40		NOT FOUND							
41		NOT FOUND							

MBS case 3969 3/6/85

V-60

279

MASS CHROMATOGRAMS

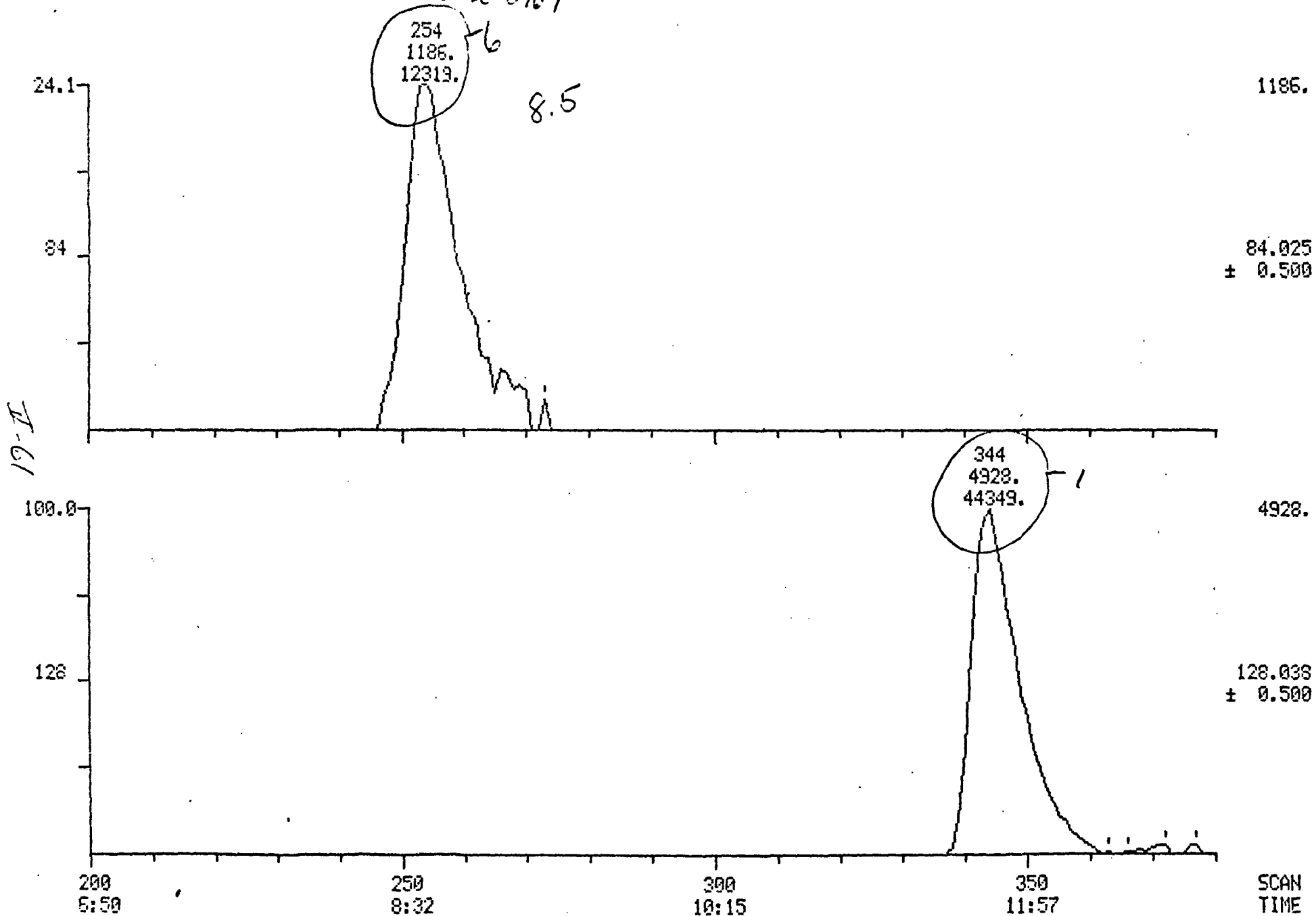
03/05/85 10:53:00

SAMPLE: BLANK 3/6/85

Case 3969

DATA: U2754

SCANS 200 TO 380



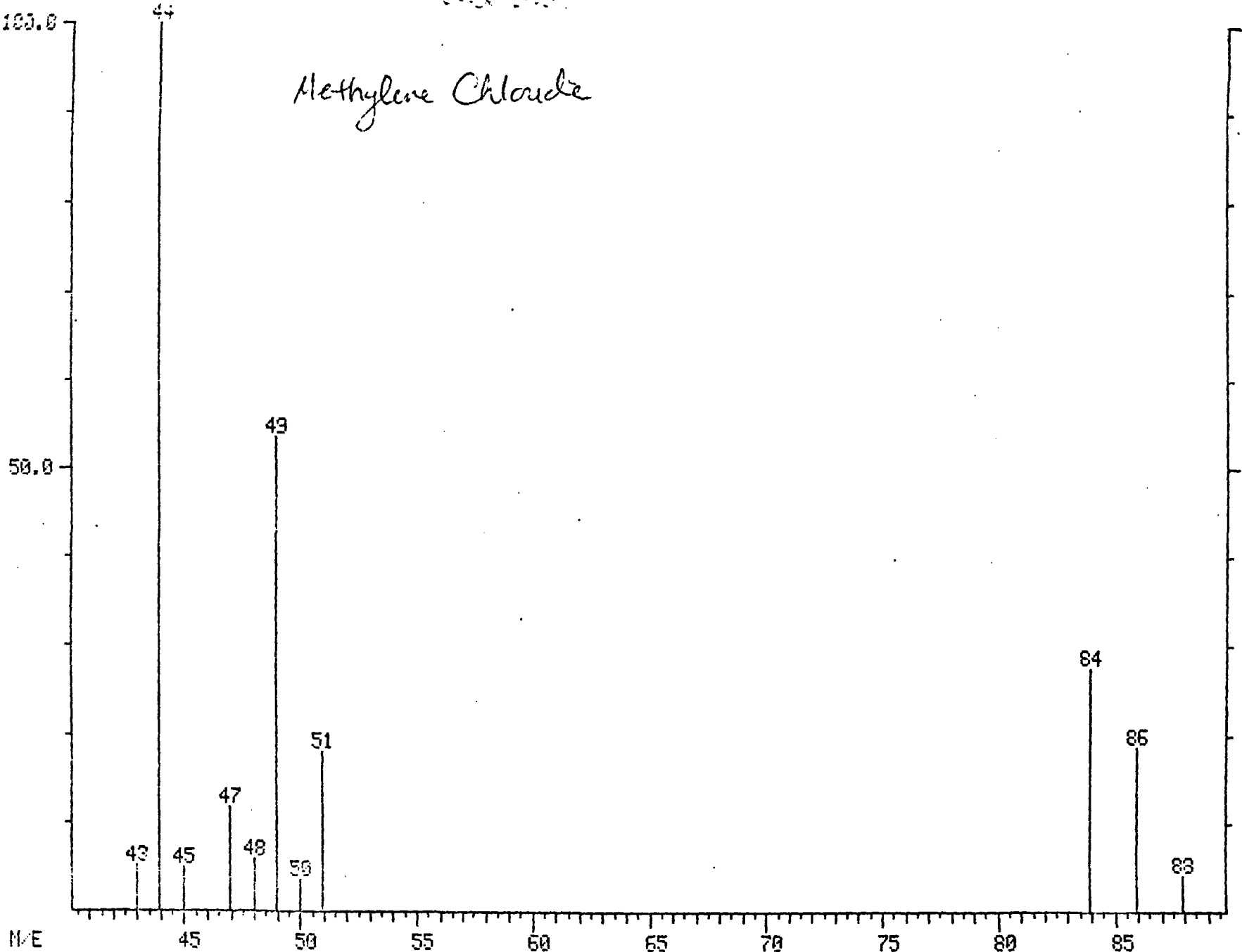
MASS SPECTRUM
08:05:35 10:50:20
SAMPLE: 11-10-85

DATE: 11-10-85

TIME: 10:50:20

27-2

Methylene Chloride



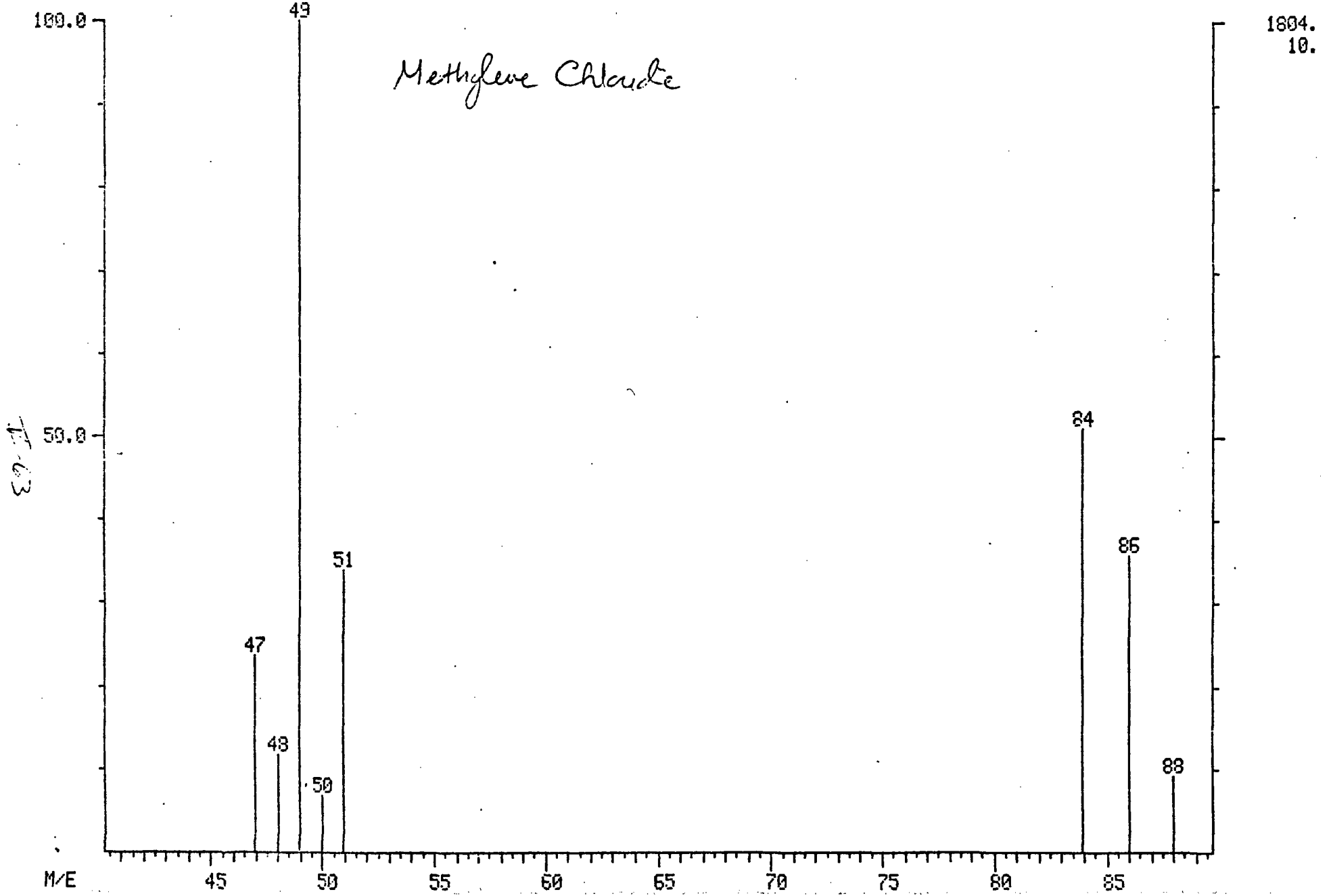
4320.
10.

MASS SPECTRUM
03/05/85 10:53:00 + 8:41
SAMPLE: SLANK 3/6/85 Case 3969
ENHANCED (S 158 2N)

DATA: U2754 #254

BASE M/E: 49
RIC: 4920.

Methylene Chloride



49.03

Handwritten mark

MASS SPECTRUM

03/03/85 12:41:00 + 8:28

SAMPLE:

ENHANCED (S 15B 2N)

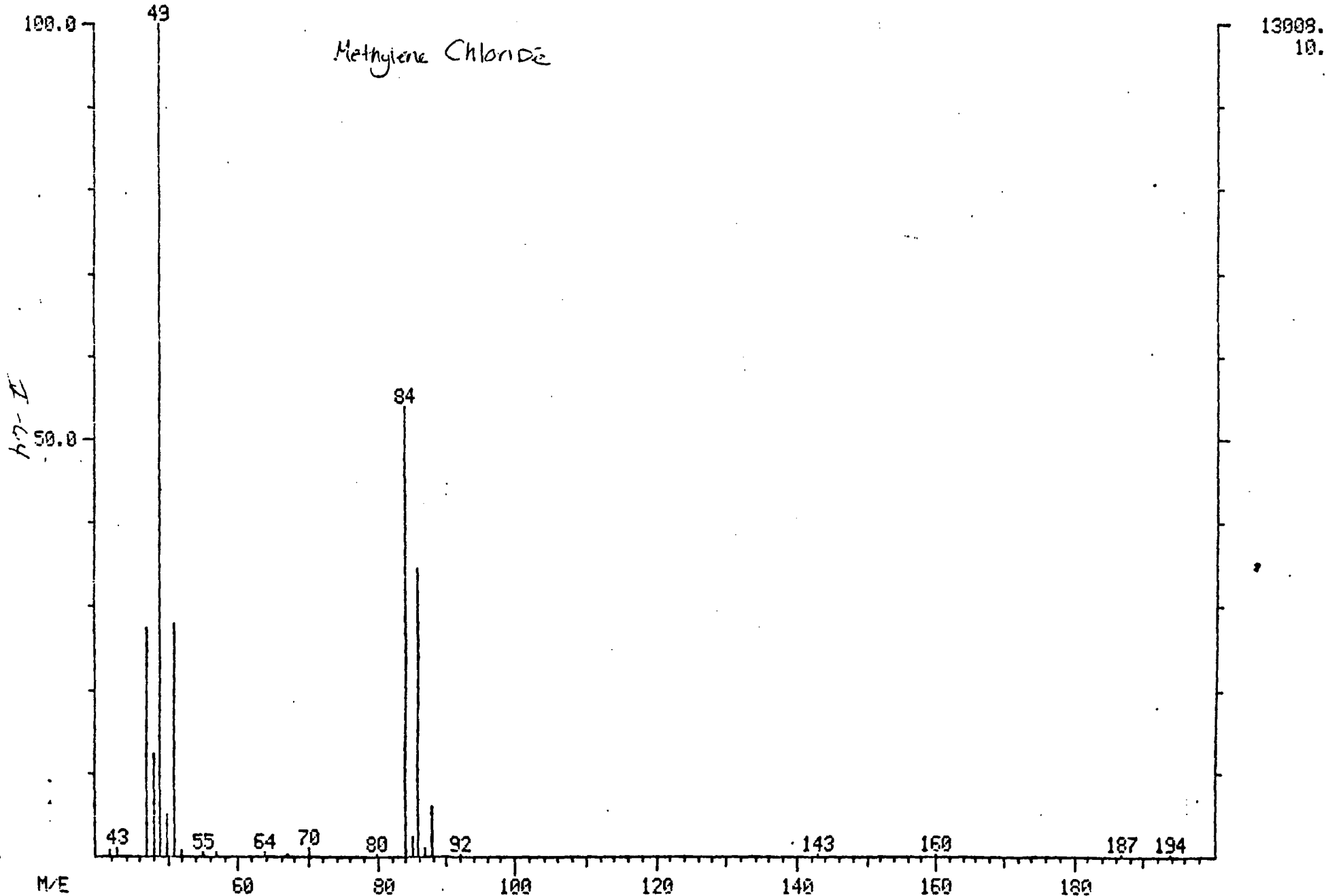
396.9
COC ~~35~~ 100 ppb

DATA: U2700 #248

BASE M/E: 49

RIC: 36224.

Methylene Chloride



h.v. - I

W

Client Name: GCA
File ID No: 433 32 MS
Matrix: Water
Date Authorized By: _____

Case No: 3969
QC Report No: 3969
Contract No: 68-01-6767
Date Sample Received: 3/1/85

Volatil Compounds

Concentration: Low Medium (Circle One)
Date Extracted/Prepared: 3/5/85
Date Analyzed: 3/5/85
Conc/Dil Factor: 1 pH: NA
Percent Moisture: NA
Percent Moisture (Determined): NA

CAS Number
1-B
2-B
3-C
4-D
5-E
6-F
7-G
8-H
9-I
10-J
11-K
12-L
13-M
14-N
15-O
16-P
17-Q
18-R
19-S
20-T
21-U
22-V
23-W
24-X
25-Y
26-Z

CAS Number	Compound	ug/l or ug/Kg (Circle One)
1-B	Chloromethane	10.1
2-B	Bromomethane	10.1
3-C	Vinyl Chloride	10.1
4-D	Chloroethane	10.1
5-E	Methylene Chloride	4.5
6-F	Acetone	11.1
7-G	Carbon Disulfide	5.1
8-H	1,1-Dichloroethane	5.1
9-I	1,1-Dichloroethane	5.1
10-J	Trans-1,2-Dichloroethane	5.1
11-K	Chloroform	6.8
12-L	1,2-Dichloroethane	5.1
13-M	2-Butanone	10.1
14-N	1,1,1-Trichloroethane	5.1
15-O	Carbon Tetrachloride	5.1
16-P	Vinyl Acetate	10.1
17-Q	Bromodichloromethane	5.1

CAS Number	Compound	ug/l or ug/Kg (Circle One)
79-34-5	1,1,2,2-Tetrachloroethane	5.1
78-87-5	1,2-Dichloropropane	5.1
10061-09-6	Trans-1,3-Dichloropropene	5.1
79-01-5	Trichloroethylene	5.1
124-48-1	Dibromochloromethane	5.1
79-00-5	1,1,2-Trichloroethane	5.1
71-43-2	Benzene	5.1
10051-01-5	cis-1,3-Dichloropropene	5.1
110-25-8	2-Chloroethylvinyl ether	10.1
75-25-2	Bromoforn	5.1
591-76-6	2-Hexanone	5.1
108-10-1	4-Methyl-2-Pentanone	5.1
117-18-4	Tetrachloroethene	5.1
108-88-3	Toluene	5.1
108-89-7	Chlorobenzene	5.1
100-41-4	Ethylbenzene	5.1
100-42-5	Styrene	5.1
	Total Xylenes	5.1

For a complete explanation of the results, please refer to the report and the laboratory manual.

- 1. Indicates a value greater than the detection limit reported for this compound.
- 2. Indicates a compound was detected but not identified. Report the compound name and the detection limit (e.g., 100 ug/l based on mass) and the concentration of the compound (e.g., 10 ug/l based on mass) in the instrument detection limit. The detection limit for this compound was analyzed. If the compound was detected, the number in the instrument detection limit for the sample.
- 3. Indicates an estimated value. This value is used when the compound is not identified but the mass is known. It is used when the mass is known but the compound is not identified. The number in the instrument detection limit for the sample.
- 4. This flag is used to indicate parameters where the detection limit has been confirmed by GC-MS. Single sample parameters for 10 ug/l in the final extract should be confirmed by GC-MS.
- 5. This flag is used when the analysis found in the final extract is not identified. It indicates possible presence but is not confirmed. It warns the data user to take a proper action.
- 6. Other specific comments or footnotes may be required to provide details on the results. If used, they must be fully documented in the report.

IT-65

Sample Number
AB 028.ms

Organics Analysis Data Sheet
 (Page 2)

Semivolatiles Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 3/5/95

Date Analyzed: 3/25/95

Conc/Dil Factor: 500mls -> 1.0ml

CA# Number	Compound Name	ug/l or ug/Kg (Circle One)
1	N-Nitrosodimethylamine	10u
2	Ethanol	
3	Aniline	10u
4	tert-2-Chloroethyl Ether	10u
5	2-Chlorophenol	
6	1,3-Dichlorobenzene	10u
7	1,4-Dichlorobenzene	
8	Benzyl Alcohol	10u
9	1,2-Trichlorobenzene	10u
10	2-Methylphenol	10u
11	Bis(2-chloro-propyl) Ether	10u
12	4-Methylphenol	10u
13	N-Nitroso-Di-n-Propylamine	
14	Hexachlorocyclopentadiene	10u
15	1,2,3-Trichlorobenzene	10u
16	Isophorone	10u
17	2-Nitrophenol	10u
18	2,4-Dimethylphenol	10u
19	Benzoic Acid	50u
20	tert-2-Chloroethyl Methane	10u
21	2,4-Dichlorophenol	10u
22	1,2,4-Trichlorobenzene	
23	Naphthalene	10u
24	4-Chloroaniline	10u
25	Hexachlorobutadiene	10u
26	4-Chloro-3-methylphenol	
27	2-Methylanthralene	10u
28	Hexachlorocyclopentadiene	10u
29	2,4,6-Trichlorophenol	10u
30	2,4,5-Trichlorophenol	50u
31	1-Chloroanthracene	10u
32	2-Nitroaniline	50u
33	Dimethyl Phthalate	10u
34	Acenaphthylene	10u
35	3-Nitroaniline	50u

CAS Number	Compound Name	ug/l or ug/Kg (Circle One)
83-32-9	Acenaphthone	
81-28-5	2,4-Dinitrophenol	50u
100-02-7	4-Nitrophenol	
132-64-9	1,3-Benzofuran	10u
121-14-2	2,4-Dinitrotoluene	
806-20-2	2,6-Dinitrotoluene	10u
84-66-2	Diethylphthalate	10u
7005-72-3	4-Chlorophenyl-phenylether	10u
86-73-7	Fluorene	10u
100-01-6	4-Nitroaniline	50u
534-52-1	4,6-Dinitro-2-methylphenol	10u
86-80-6	N-Nitrosodiphenylamine (1)	10u
101-55-3	4-Propoxyphenyl-phenylether	10u
118-74-1	Hexachlorocyclopentadiene	10u
87-86-5	Pentachlorophenol	
85-01-8	Phenanthrene	10u
120-12-7	Anthracene	10u
84-74-2	Di-n-Butylphthalate	
205-44-0	Fluoranthene	10u
82-87-5	Benzidine	50u
129-00-0	Pyrene	
85-68-7	Benzylbenzylphthalate	10u
31-94-1	3,3'-Dichlorobenzidine	50u
86-55-3	Benzofluoranthene	10u
117-81-7	Bis(2-Ethylhexyl)phthalate	10u
218-01-9	Chrysene	10u
117-84-0	Di-n-Octyl Phthalate	10u
205-99-2	Benzofluoranthene	10u
207-08-9	Benzofluoranthene	10u
10-32-8	Benzofluoranthene	10u
193-39-5	Indeno(1,2,3-cd)Pyrene	10u
83-70-3	Dibenz(a,h)Anthracene	10u
191-24-2	Benzofluoranthene	10u

(1) Cannot be separated from diphenylamine

Sample Number
AB030MS

Organics Analysis Data Sheet
 (Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 3-1-85

Date Analyzed: 4/2/85

Conc/Dil Factor: 1

CAS Number ug/L or ug/Kg
(Circle One)

319-84-6	Alpha-BHC	0.050
319-85-7	Beta-BHC	0.050
319-86-8	Delta-BHC	0.050
58-69-9	Gamma-BHC (Lindane)	0.050
76-44-8	Heptachlor	0.050
509-90-2	Aldrin	0.050
1024-57-3	Heptachlor Epoxide	0.050
959-98-8	Endosulfan I	0.050
50-57-1	Dieldrin	0.050
72-85-9	4,4'-DDE	0.050
72-20-8	Endrin	1.00
39213-65-8	Endosulfan II	0.050
72-54-8	4,4'-DDD	0.050
7421-93-4	Endrin Aisobutyl	0.050
1031-07-8	Endosulfan Sulfate	0.050
50-29-3	4,4'-DDT	1.00
72-43-5	Methoxychlor	0.050
58494-70-5	Endrin Keptone	0.050
57-74-9	Chlordane	0.050
8001-35-7	Toxaphene	1.00
12674-11-2	Aroclor-1016	0.050
11104-28-2	Aroclor-1216	0.050
11141-10-5	Aroclor-1232	0.050
53459-21-0	Aroclor-1248	0.050
12672-29-6	Aroclor-1254	0.050
11097-69-1	Aroclor-1260	1.00
11095-82-5	Aroclor-1268	1.00

J. H. L. L. L.
 4/3/85



V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

V_s 500 ml BL or W_s 1000 V_i 10,000 ul V_t 5.0 ul

II 27

Organics Analysis Data Sheet
(Page 4)

AB 008 MS

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	unknown	VOA	824	31
2.				
3.				
4.				
5.				
6.				
7.				
8.				
9.				
10.				
11.				
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17.				
18.				
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20.				
21.				
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23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				

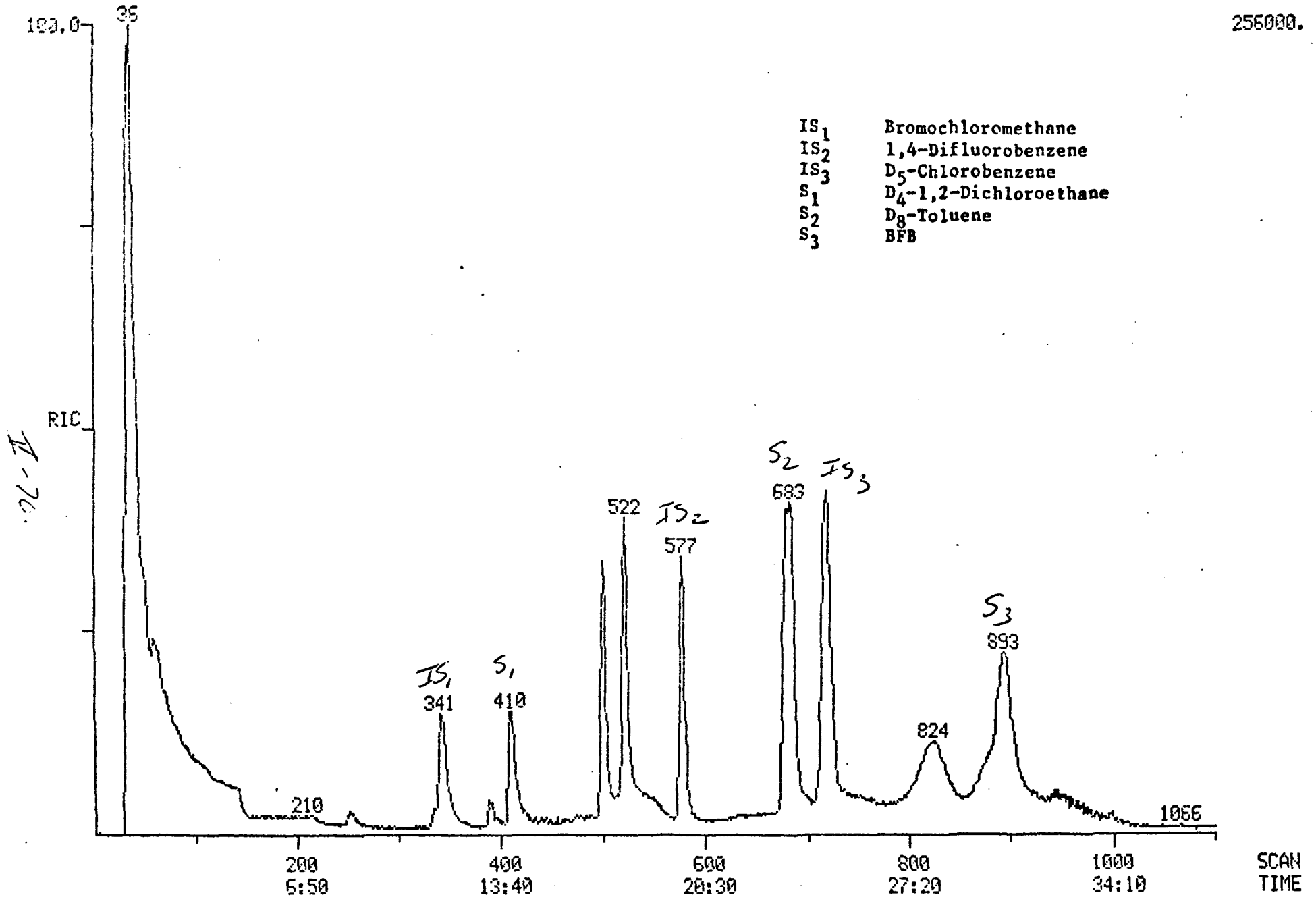
II-68

RIC
03/05/85 14:00:00
SAMPLE: 43232 MS CASE 3969 AB-028 MS

DATA: U2741

SCANS 1 TO 1100

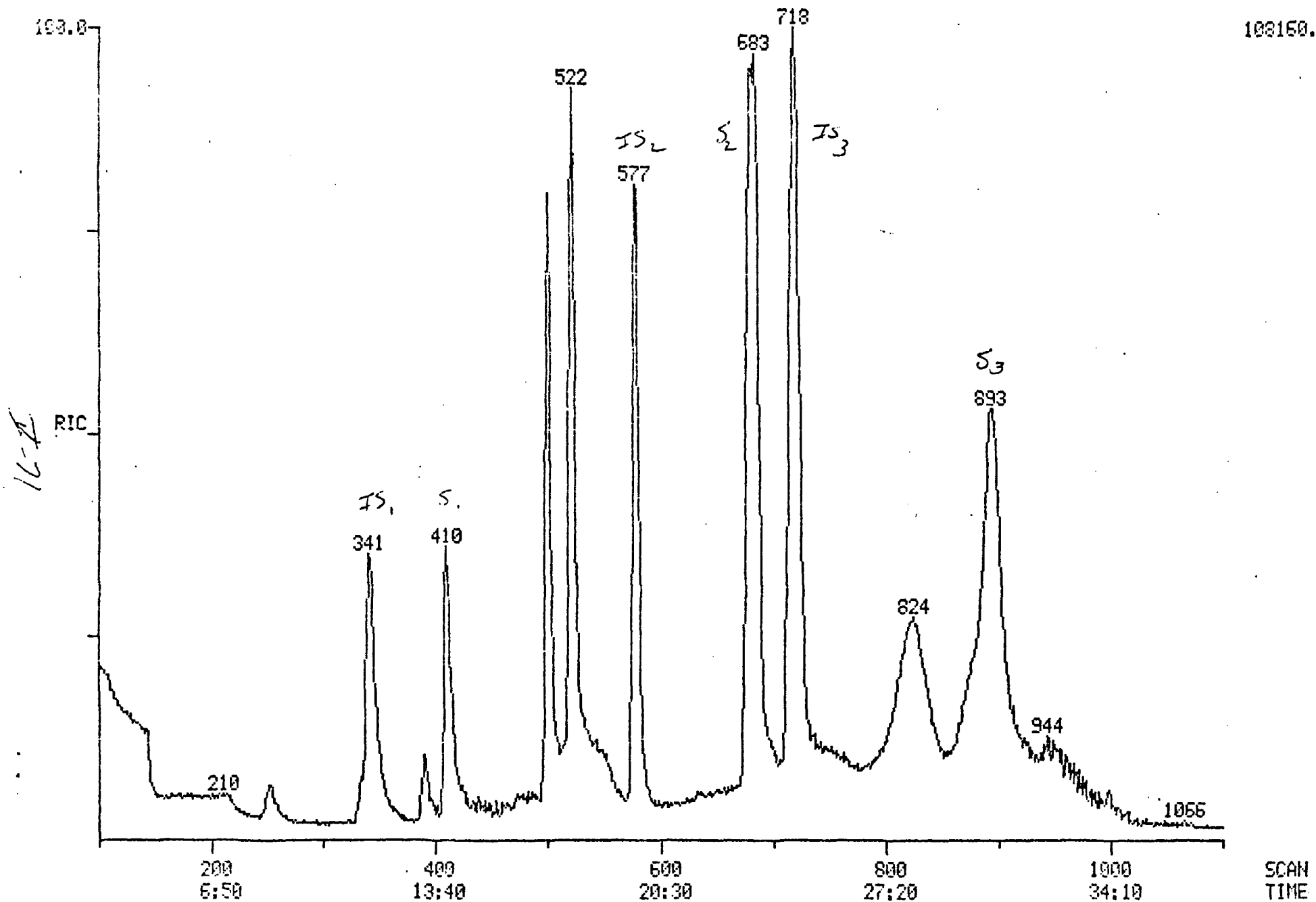
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RIC
03/05/85 14:00:00
SAMPLE: 43232 MS CASE 3359 AB-028 MS

DATA: U2741

SCANS 100 TO 1100



DATA: V2741.TI
 03/05/85 14:00:00
 SAMPLE: 43232 MS CASE 3969 AB-028 MS
 SUBMITTED BY: ANALYST:

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

- NO NAME
- 1 BROMOCHLOROMETHANE (I. S. #1)
- 2 CHLOROMETHANE
- 3 BROMOMETHANE
- 4 VINYL CHLORIDE
- 5 CHLOROETHANE
- 6 METHYLENE CHLORIDE
- 7 ACETONE
- 8 CARBON DISULFIDE
- 9 1, 1-DICHLOROETHENE
- 10 1, 1-DICHLOROETHANE
- 11 TRANS 1, 2-DICHLOROETHENE
- 12 CHLOROFORM
- 13 1, 2-DICHLOROETHANE
- 14 D4-1, 2-DICHLOROETHANE
- 15 1, 1, 1-TRICHLOROETHANE
- 16 1, 4-DIFLUOROBENZENE (I. S. #2)
- 17 CARBON TETRACHLORIDE
- 18 VINYL ACETATE
- 19 2-BUTANONE
- 20 BROMODICHLOROMETHANE
- 21 1, 2-DICHLOROPROPANE
- 22 TRANS 1, 3-DICHLOROPROPANE
- 23 TRICHLOROETHENE
- 24 DIBROMOCHLOROMETHANE
- 25 1, 1, 2-TRICHLOROETHANE
- 26 BENZENE
- 27 CIS-1, 3-DICHLOROPROPENE
- 28 BROMOFORM
- 29 D5-TOLUENE (SURR#2)
- 30 2-CHLOROETHYL VINYLETHER
- 31 D5-CHLOROBENZENE (I. S. #3)
- 32 4 METHYL 2 PENTANONE
- 33 2-HEXANONE
- 34 TETRACHLOROETHENE
- 35 1, 1, 2, 2, TETRACHLOROETHANE
- 36 TOLUENE
- 37 CHLOROBENZENE
- 38 EFB (SURR#3)
- 39 ETHYLBENZENE
- 40 STYRENE
- 41 TOTAL XYLENES

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	125	341	11:39	1	1.000	A BB	30310.	50.000 UC/L	7.71
2									
3									

V-72

NO	SCAN	TIME	REF	RRT	METH	AREA (HGHT)	AMOUNT	%TOT
NO	OUND							
NO	OUND							
NO	OUND							
NO	OUND							
NO	OUND							
NO	OUND							
NO	OUND							
8	OUND							
NO	390	13:19	1	1.144	A BB	21691.	6.772 UG/L	1.04
6	OUND							
NO	411	14:03	1	1.205	A BB	105209.	95.739 %REC	14.76
114	OUND							
NO	577	19:43	16	1.000	A BB	151476.	50.000 UG/L	7.71
NO	OUND							
NO	OUND							
NO	OUND							
NO	OUND							
130	OUND							
NO	500	17:05	16	0.867	A BV	61214.	52.429 UG/L	8.08
NO	OUND							
7	OUND							
NO	522	17:50	16	0.905	A BB	189007.	45.634 UG/L	7.04
NO	OUND							
9	OUND							
NO	679	23:12	16	1.177	A BB	176993.	96.100 %REC	14.82
117	OUND							
NO	716	24:28	31	1.000	A BB	110130.	50.000 UG/L	7.71
NO	OUND							
5	OUND							
164	678	23:10	31	0.947	A BB	2723.	6.081 UG/L	0.94
8	634	21:40	31	0.885	A BB	126.	0.127 UG/L	0.02
9	639	21:50	31	0.892	A BB	122.	0.115 UG/L	0.02
11	684	23:22	31	0.955	A BB	120980.	48.197 UG/L	7.43
9	720	24:36	31	1.006	A BB	126354.	50.555 UG/L	7.79
10	854	30:33	31	1.249	A BB	93725.	95.280 %REC	14.69
NO	809	27:38	31	1.130	A BV	2023.	1.588 UG/L	0.24
NO	OUND							
NO	OUND							

AB 028 MS Case 3969 3/5/85

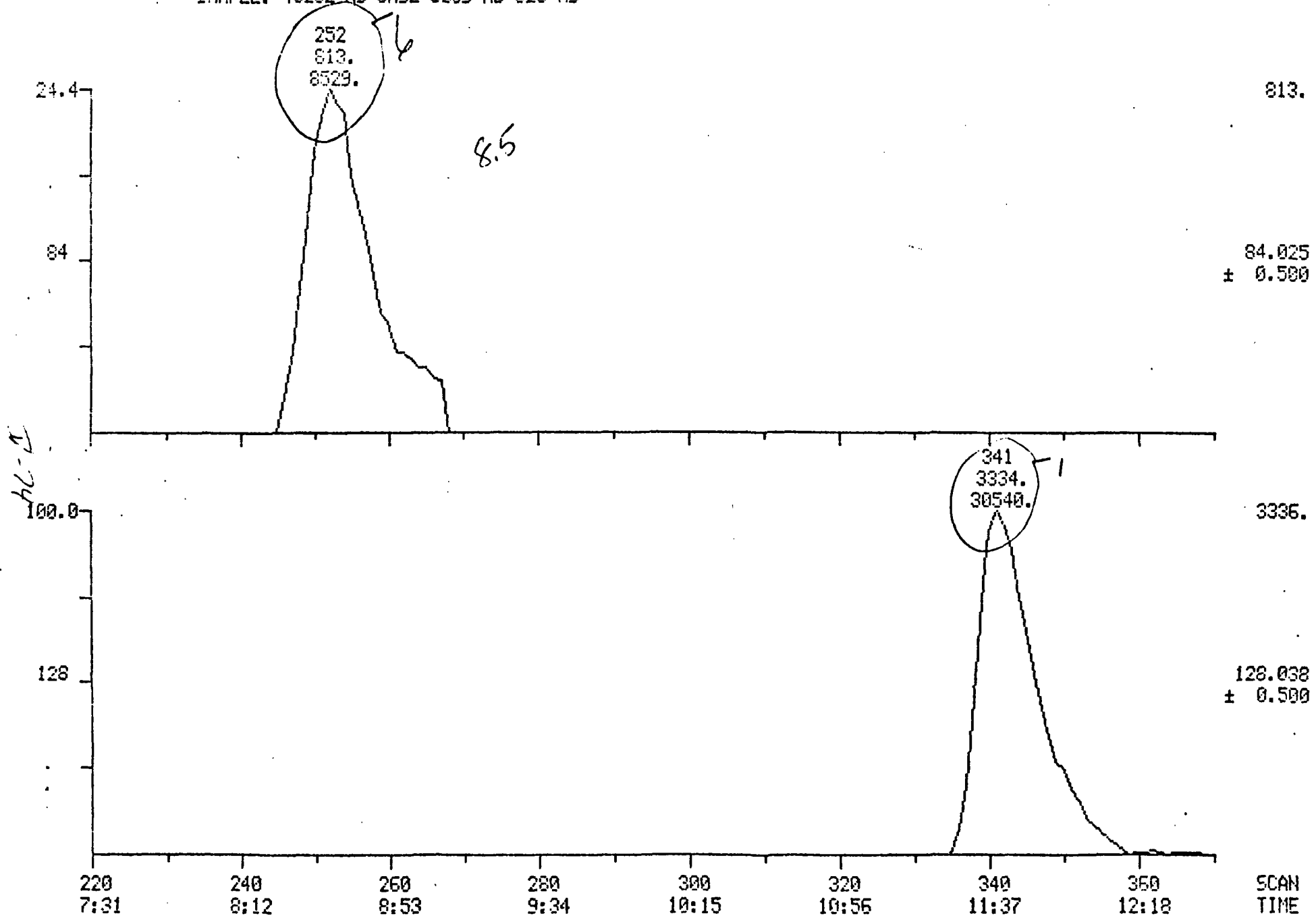
V-73

257

MASS CHROMATOGRAMS
03/05/85 14:00:00
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DATA: U2741

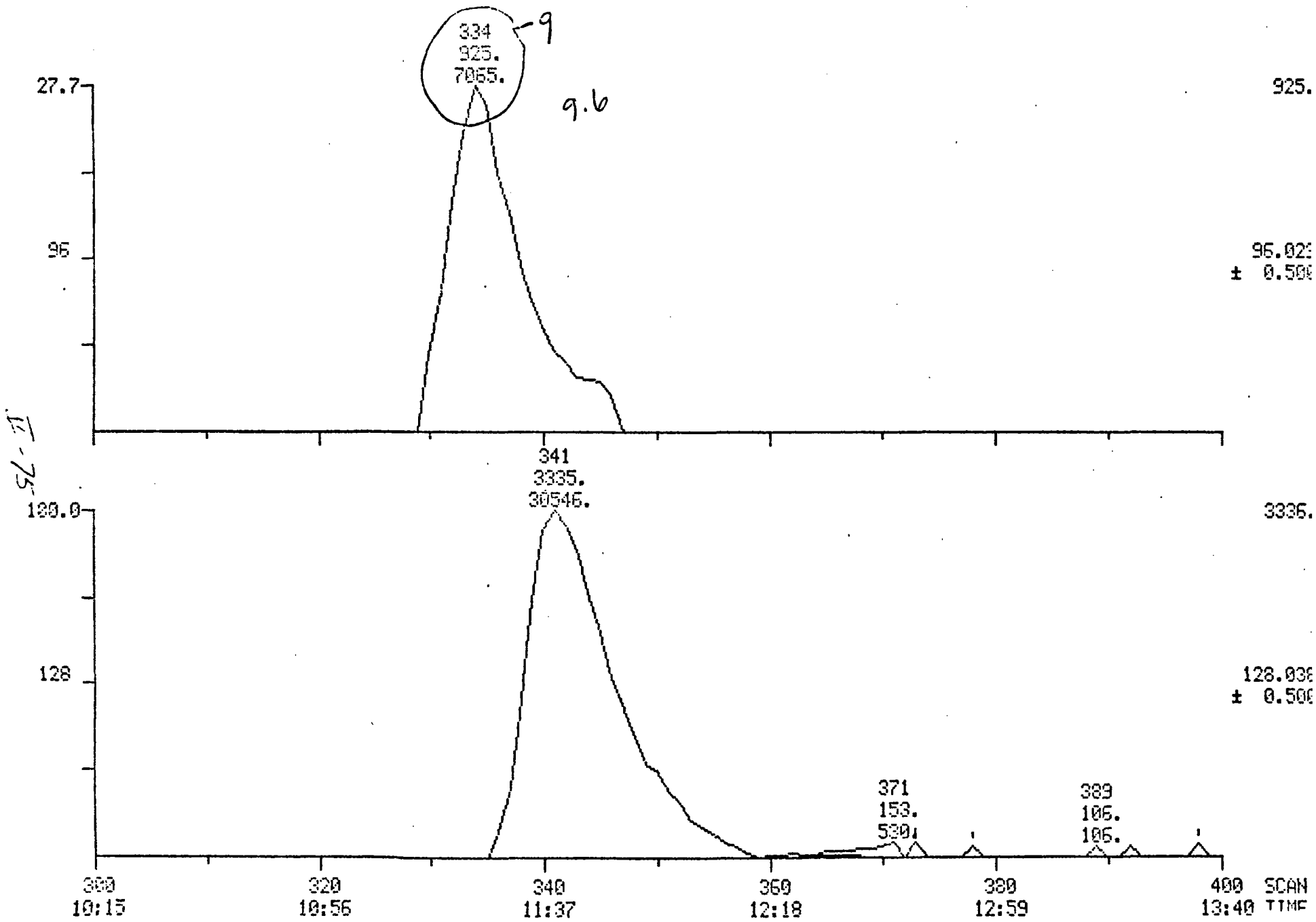
SCANS 220 TO 370



MASS CHROMATOGRAMS
03/05/85 14:00:00
SAMPLE: 43232 MS CASE 3959 AB-028 MS

DATA: U2741

SCANS 300 TO 400



LIBRARY SEARCH
03/05/85 14:00:00 + 28:09
SAMPLE: 43232 MS CASE 3969 AB-028 MS
ENHANCED (S 158 2N 0T)

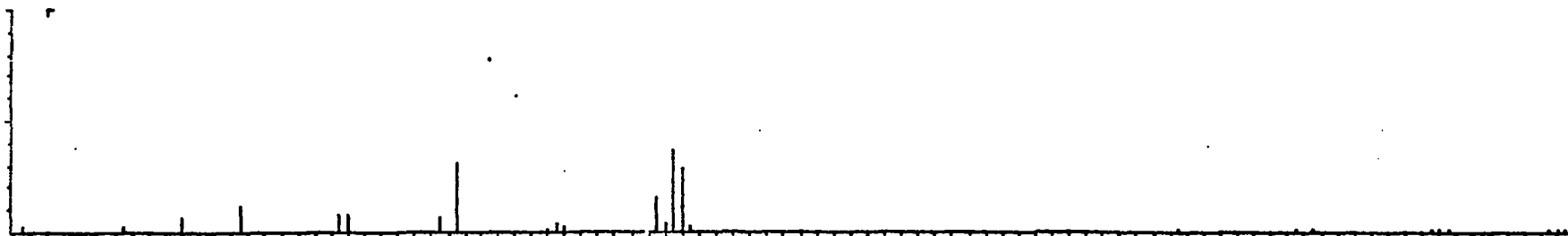
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BASE M/E: 105
RIC: 2375.



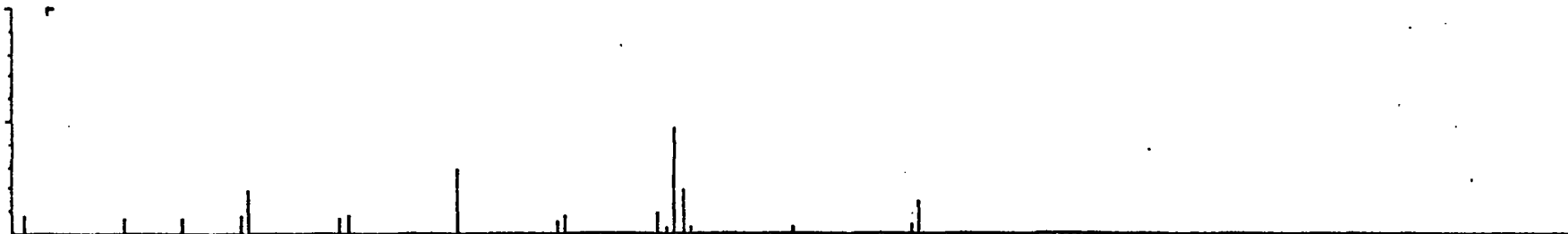
C17.H18
M WT 122
B PK 117
RANK 1
IN 15357
FIT 651

BENZENE, 1,1'-(1-ETHENYL-1,3-PR



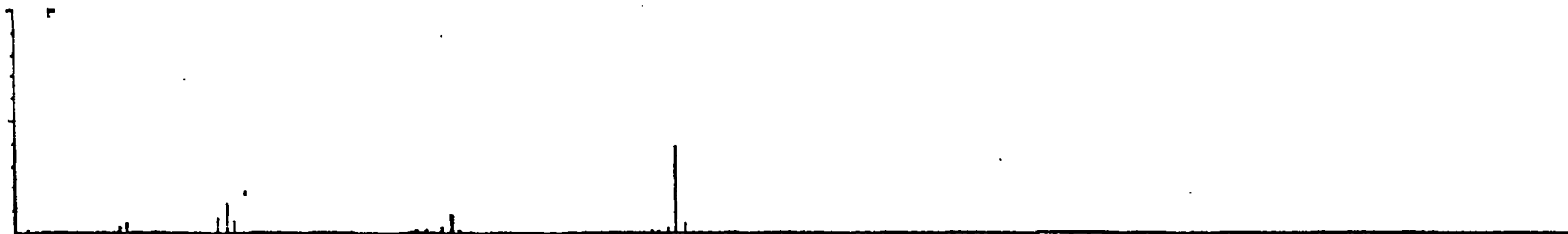
C10.H10.0
M WT 140
B PK 117
RANK 2
IN 5416
FIT 624

1,2,4-METHENO-3H-CYCLOBUTA[CD]



C8.H7.N
M WT 117
B PK 117
RANK 3
IN 2300
FIT 588

1H-INDOLE



M/E 40 60 80 100 120 140 160 180 200 220

Handwritten: 76 77 78

Handwritten: 200

03/05/85 14:00:00
SAMPLE: 43232 MS CASE 3959 AB-028 MS

DATA: 02741

SCAN: 000 TO 000

825
3128.
67641.

$$\frac{67641}{110130} \times \frac{50}{1} \approx 31$$

3128.

115

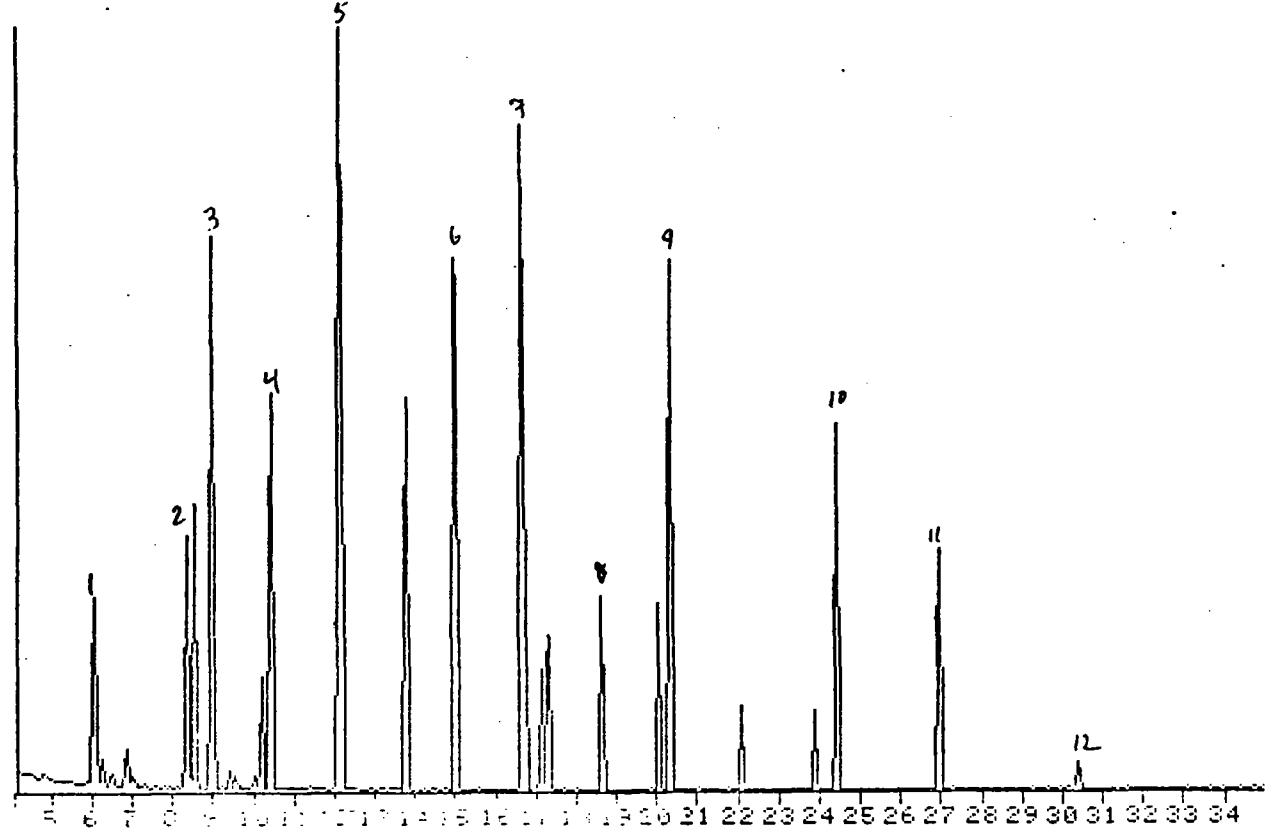
100.0
185

105.031
± 0.500

800 27:20 810 27:40 820 28:01 830 28:21 840 28:42 850 29:02 860 29:23 SCAN TIME

83847

ST-16



1. 2-FLUOROPHENOL (SURF)
2. PHENOL-D6 (SURR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SURF)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUOROBIPHENYL (SURF)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SURR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SURF)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

V-78

✓

FILE NUMBER 14757
 QC-957 RB028MS CASE 3969
 3/25/85/EM BTL#16 Q14757 D14757

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
9.0	150.0	1.000	80.0000				D4-1,4-DICHLOROBENZENE(20)
9.9	74.0	.046	.0000				N-NITROSODIMETHYLAMINE
8.4	93.0	.913	.0000				ANILINE
8.5	93.0	.709	1.8523 EM				BIS(2-CHLOROETHYL)ETHER
8.9	146.0	.591	.0000				1,3-DICHLOROBENZENE
9.0	146.0	.594	46.9800 ✓				1,4-DICHLOROBENZENE
9.5	108.0	.395	.0000				BENZYL ALCOHOL
9.5	146.0	.594	.0000				1,2-DICHLOROBENZENE
9.9	45.0	.527	.0000				BIS(2-CHLOROISOPROPYL)ETHER
10.2	70.0	.490	55.3977 ✓				N-NITROSODINPROPYLAMINE
10.2	117.0	.291	.0000				HEXACHLOROETHANE
10.4	77.0	.716	.5272				NITROBENZENE

MORE ?

12.1	68.0	1.000	80.0000				D8-NAPHTHALENE(20)
11.1	138.0	1.519	.0000				ISOPHORONE
11.8	95.0	1.813	.0000				BIS(2-CHLOROETHOXY)METHANE
12.1	180.0	2.642	42.4707 ✓				1,2,4-TRICHLOROBENZENE
12.2	129.0	1.325	.0000				NAPHTHALENE
12.5	127.0	4.577	.0000				4-CHLOROANILINE
12.7	225.0	1.190	.0000				HEXACHLOROBUTADIENE
14.0	115.0	2.578	.0000				2-METHYLNAPHTHALENE

16.7	164.0	1.000	80.0000				D10-ACENAPHTHENE
14.6	237.0	.205	.0000				HEXACHLOROCYCLOPENTADIENE
15.2	164.0	.281	.0000				2-CHLORONAPHTHALENE
15.6	138.0	.397	.0000				2-NITROANILINE
16.2	164.0	.083	.0000				DIMETHYL PHTHALATE
16.3	151.0	.272	.0000				ACENAPHTHYLENE
16.4	165.0	.228	.0000				2,6-DINITROTOLUENE
16.7	138.0	.293	.0000				3-NITROANILINE
16.7	152.0	.437	48.0670 ✓				ACENAPHTHENE
17.2	139.0	.490	22.8977 EM				DIBENZOFURAN
17.3	165.0	.320	44.3927 ✓				2,4-DINITROTOLUENE
18.0	177.0	.171	.0000				DIETHYL PHTHALATE
18.0	165.0	.820	.0000				FLUORENE
18.1	204.0	.358	.0000				4-CHLOROPHENYLPHENYL ETHER
18.2	138.0	.292	.0000				4-NITROANILINE
17.2	169.0	.144	.0000				DIPHENYLAMINE

D-79

113

20.3	188.0	1.000	80.0000	D10-PHENANTHRENE(20)
19.3	248.0	.112	.0000	4-BROMOPHENYLPHENYL ETHER
19.6	284.0	.112	.0000	HEXACHLOROBENZENE
20.4	179.0	.060	.0000	PHENANTHRENE/ANTHRACENE
22.1	150.0	.100	14.1619✓	DIBUTYL PHTHALATE
23.4	202.0	.665	.0000	FLUORANTHENE

FILE NUMBER 14757
 QC-957 AB028MS CASE 3969
 3/25/85/EM BTL#16 Q14757 D14757

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
27.0	240.0	1.000	80.0000				D12-CHRYSENE
23.9	202.0	2.264	14.8597✓				PYRENE
25.8	149.0	1.136	.0000				BUTYL BENZYL PHTHALATE
26.9	228.0	.427	.0000				BENZO(A)ANTHRACENE/CHRYSENE
27.3	149.0	1.128	.0000				BIS(2-ETHYLHEXYL)PHTHALATE

30.4	264.0	1.000	80.0000				D12-PERYLENE
28.7	149.0	5.974	.0000				DIOCTYL PHTHALATE
29.5	252.0	1.307	.0000				BENZO(B)/(K)FLUORANTHENE
30.2	252.0	.744	.0000				BENZO(A)PYRENE
33.7	276.0	.485	.0000				INDENO(123-CD)PYRENE
33.8	278.0	.346	.0000				DIBENZO(AH)ANTHRACENE
34.6	276.0	.360	.0000				BENZO(GHI)PERYLENE

27.0	240.0	1.000	80.0000				D12-CHRYSENE
23.8	184.0	.712	.0000				BENZIDINE
27.0	252.0	.146	.0000				DICHLOROBENZIDINE

#-80

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
9.0	150.0	1.000	80.0000				D4-1,4-DICHLOROBENZENE(20)
8.4	94.0	.841	56.7688	✓			PHENOL
8.5	128.0	.611	87.9009	✓			2-CHLOROPHENOL
9.9	108.0	.671	.0000				2-METHYLPHENOL
10.2	108.0	.699	.0000				4-METHYLPHENOL

12.1	68.0	1.000	80.0000				D8-NAPHTHALENE(20)
11.3	139.0	1.759	.0000				2-NITROPHENOL
11.5	122.0	2.581	.0000				2,4-DIMETHYLPHENOL
12.1	122.0	1.200	.0000				BENZOIC ACID
11.9	162.0	2.051	.0000				2,4-DICHLOROPHENOL
13.8	142.0	2.196	86.9397	✓			4-CHLORO-M-CRESOL

16.7	164.0	1.000	80.0000				D10-ACENAPHTHENE(20)
14.8	196.0	.237	.0000				2,4,5-TRICHLOROPHENOL
14.8	196.0	.237	.0000				2,4,6-TRICHLOROPHENOL
17.0	184.0	.111	.0000				2,4-DINITROPHENOL
18.3	198.0	.126	.0000				4,6-DINITRO-O-CRESOL
17.2	65.0	.231	49.1382	✓			4-NITROPHENOL

20.3	188.0	1.000	80.0000				D10-PHENANTHRENE
20.1	266.0	.072	81.2828	✓			PENTACHLOROPHENOL

Date 4/1/75Analyst BLContract 3-705-050

Pesticide Analysis - Qualitative/Quantitative

Case # 3969Sample I.D. CC-967/AB030 MS
Run # 15HP 5870
Bottle # 15

RT (min.)	Peak ID	Area $\times 10^3$	ng/ml	Dilution Factor (mls extracted)	Spiked at μg	% ($\mu\text{g}/\text{l}$) ppb Recovery	total μg
2.58	δ -BHC	26.50	* $12.8 \times 10 = 128$		200	64.0	
3.18	Heptachlor	34.07	$13.0 \times 10 = 130$		200	65.0	
3.87	Aldrin	31.68	$15.5 \times 10 = 155$		200	77.5	
9.42	Dieldrin	64.09	$39.0 \times 10 = 390$		500	78.0	
11.52	Endrin	63.80	$54.0 \times 10 = 540$		500	107	
17.34	pp'-DDT	41.83	* $55.2 \times 10 = 552$		500	110	
31.70	DBC	203.5	$\circ 154 \times 10 = 1540$		1000	154	
* Calculated from lowest four calibration points							
* Calculated from highest four calibration points							
\circ There is a shoulder on the DBC peak at 31.70 min which gave the high area count and corresponding high recovery.							

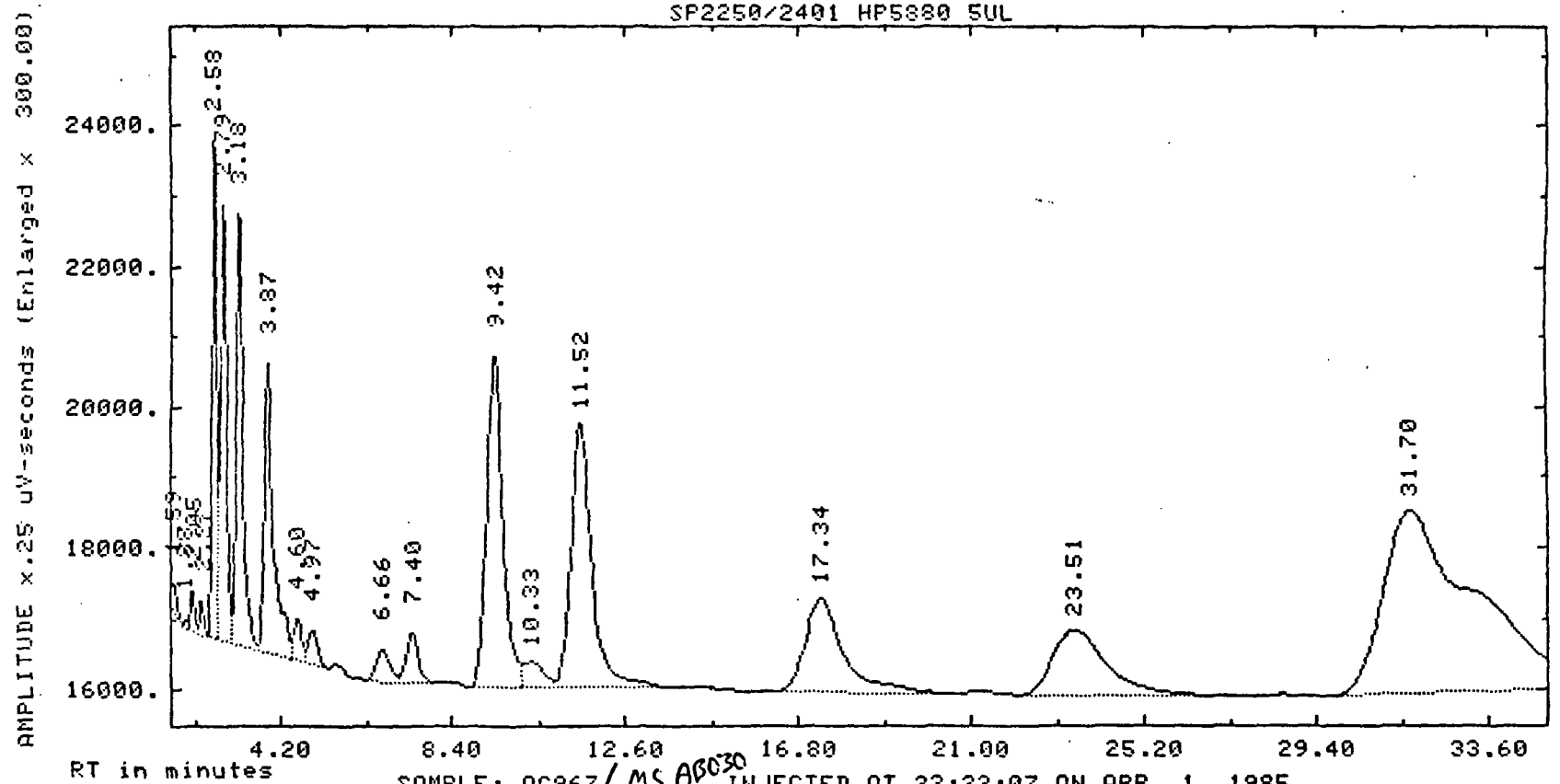
GCA
GCAGCA CORPORATION
Technology Division

J 92

EB-II

Case 3969

SP2250/2401 HP5880 5UL



SAMPLE: QC967/MS A8030 INJECTED AT 23:22:07 ON APR 1, 1985
Method: PRIME Raw: RB5015 Proc: PB5015

*LI,P,PB5015

PROCESSED DATA FILE: PB5015 ON CRN 21

APR 2, 1985 9:13:47

Case 3969

REPORT: 150 CHANNEL: 2 # PEAKS: -21 SP2250/2401 HP5880 SUL

PEAK	RT	ITM	FACTOR	AREA	AREA %	NAME
1	1.09	0.00	1.00000	3475 BV	.607	
2	1.33	0.00	1.00000	668 VV	.117	
3	1.40	0.00	1.00000	966 VV	.168	
4	1.50	0.00	1.00000	1709 VV	.312	
5	1.07	0.00	1.00000	292 VV	.049	
6	2.05	0.00	1.00000	2076 VV	.362	
7	2.27	0.00	1.00000	1572 VV	.274	
8	2.50	0.00	1.00000	26504 VV	4.529	<i>γ-BHC</i>
9	2.79	0.00	1.00000	30059 VV	5.251	
10	3.10	0.00	1.00000	34060 VV	5.949	<i>Heptachlor</i>
11	3.07	0.00	1.00000	31676 VV	5.532	<i>Aldrin</i>
12	4.60	0.00	1.00000	3626 VV	.658	
13	4.97	0.00	1.00000	3744 VB	.654	
14	6.60	0.00	1.00000	5282 BV	.919	
15	7.40	0.00	1.00000	7124 VB	1.244	
16	9.42	0.00	1.00000	64090 BV	11.192	<i>Dieldrin</i>
17	10.33	0.00	1.00000	5435 VV	.949	
18	11.52	0.00	1.00000	63001 VB	11.142	<i>Endrin</i>
19	17.34	0.00	1.00000	41020 BB	7.305	<i>pp'-DDT</i>
20	23.51	0.00	1.00000	40901 BB	7.136	
21	31.70	0.00	1.00000	203504 BF	35.539	<i>Dibutylchlorodate</i>

TOTAL AREA = 572620 TOTAL AREA % = 100.000

SAMPLE: Q0507 / MS A030 INJECTED AT 23:22:07 ON APR 1, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 15 GTL: 15

SL-WDTH	NU/MIN	DELAY	MIN-AR	BUNCH	REPORT
.500	.300	1.00	50	AUTO	MEDIUM

SUP-UNK	DVT	ID-LVL	REF-RTW	%RTW	%DIL-F	ISO
NO	0.00	0	.30	5.0	100.00	YES

ACTUAL RUN TIME: 40.017 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: R05015 PARAM FILES= METHOD: SEQ:

NAME
DATE

V-84

Sample Number
AB028 MSD

Organics Analysis Data Sheet
 (Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 3/5/85

Date Analyzed: 3/25/85

Conc/Dil Factor: 500 ml to 10 ml

CAS Number		ug/l or ug/Kg (Circle One)
62-75	N-Nitrosodimethylamine	100
108-9	Phenol	
62-53	Aniline	100
111-4	Di(2-Chloroethyl)Ether	100
95-57	2-Chlorophenol	
541-7	1,2-Dichlorobenzene	100
106-46	1,4-Dichlorobenzene	
100-51	Benzyl Alcohol	100
95-50	1,2-Dichlorobenzene	100
95-48	2-Methylphenol	100
39638	Di(2-chloroisopropyl)ether	100
106-43	2-Methylphenol	100
621-6	N-Nitroso-Di-n-Propylamine	
67-72	Hexachloroethane	100
98-95	Benzobenzene	100
78-59	Isophorene	100
88-75	2-Nitrophenol	100
105-67	2,4-Dimethylphenol	100
65-85	Benzoic Acid	100
111-91	Di(2-Chloroethoxy)Methane	100
120-83	2,4-Dichlorophenol	100
120-82	1,2,4-Trichlorobenzene	
91-20	Isopthalene	100
106-47	4-Chloroaniline	100
87-65	Hexachlorobutadiene	100
69-50	4-Chloro-3-Methylphenol	
91-57	2-Methylnaphthalene	100
77-47	Hexachlorocyclopentadiene	100
83-00	2,4,6-Trichlorophenol	100
95-95	2,4,5-Trichlorophenol	100
91-58	4-Chloroanthalene	100
81-74	2-Nitroaniline	100
131-11	Dimethyl Phthalate	100
208-96	Acetyl benzydene	100
89-09	5-Nitroaniline	100

CAS Number		ug/l or ug/Kg (Circle One)
83-32	Acenaphthene	
51-28	2,4-Dinitrophenol	100
100-02	4-Nitrophenol	
132-64	Dibenzofuran	100
121-14	2,4-Dinitrotoluene	
606-20	2,6-Dinitrotoluene	100
84-66	Diethylphthalate	100
7095-72	4-Chlorophenylphenylether	100
86-73	Fluorene	100
100-01	4-Nitroaniline	100
834-52	1,6-Dinitro-2-methylphenol	100
86-30	N-Nitrosodiphenylamine (1)	100
101-55	4-Bromophenylphenylether	100
118-74	Hexachlorobenzene	100
37-86	Pentachlorophenol	
95-01	Phenanthrene	100
120-12	Anthracene	100
84-74	Di-n-Butylphthalate	
206-44	Fluoranthene	100
92-87	Benidine	100
129-00	Pyrene	
85-68	Butylbenzylphthalate	100
91-94	3,3'-Dichlorodiphenylamine	100
86-85	Benzo(a)Anthracene	100
117-81	Di(2-Ethylhexyl)Phthalate	100
813-01	Chrysene	100
117-54	Di-n-Octyl Phthalate	100
205-99	Benzo(b)Fluoranthene	100
207-09	Benzo(k)Fluoranthene	100
20-32	Benzo(a)Pyrene	100
193-39	Indeno(1,2,3-cd)perylene	100
83-70	Di(2,3-h)Anthracene	100
191-24	Benzo(g,h,i)perylene	100

(1) Cannot be separated from diphenylamine

Sample Number
A8030 MSD

Organics Analysis Data Sheet
(Page 3)

Pesticide/PCBs

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: 3-1-85

Date Analyzed: 4/2/85

Conc./Dil Factor: 1

CAS Number ug/l or ug/Kg
(Circle One)

CAS Number		<u>ug/l or ug/Kg</u> (Circle One)
319-84-6	Alpha-BHC	0.050
319-05-7	Beta-BHC	0.050
319-86-8	Delta-BHC	0.050
58-89-9	Gamma-BHC (Lindane)	0.277
76-44-8	Heptachlor	0.272
309-00-2	Aldrin	0.300
1024-57-3	Heptachlor Epoxide	0.050
959-98-8	Endosulfan I	0.050
60-57-1	Dieldrin	0.277
72-55-9	4, 4'-DDE	0.100
72-26-8	Endrin	1.19
33213-65-9	Endosulfan II	0.100
72-84-8	4, 4'-DDD	0.100
7421-93-4	Endrin Aldehyde	0.100
1031-07-8	Endosulfan Sulfate	0.100
50-29-3	4, 4'-DDT	1.19
72-43-5	Methoxychlor	0.050
53494-70-5	Endrin Ketone	0.100
57-74-9	Chlordane	0.100
8001-35-2	Toxaphene	1.19
12674-11-2	Aroclor-1016	0.500
11104-28-2	Aroclor-1221	0.500
11141-16-5	Aroclor-1232	0.500
53463-21-9	Aroclor-1242	0.500
12672-29-6	Aroclor-1248	0.500
11027-69-1	Aroclor-1254	1.19
11096-82-5	Aroclor-1260	1.19

JH Hall
4/2/85

V_i = Volume of extract injected (ul)

V_s = Volume of water extracted (ml)

W_s = Weight of sample extracted (g)

V_t = Volume of total extract (ul)

500 ml or W_s V_i 10,000 ml V_t 5.0 ml

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Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	RT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	<i>No compounds found</i>		<i>NCA</i>	
2.				
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1-29

AB 028 MSD

Organics Analysis Data Sheet
(Page 4)

Tentatively Identified Compounds

CAS Number	Compound Name	Fraction	NT or Scan Number	Estimated Concentration (ug/l or ug/kg)
1.	No Compounds found	B&A		
2.				
3.				
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40.				

D-89

10/1/89

1/84

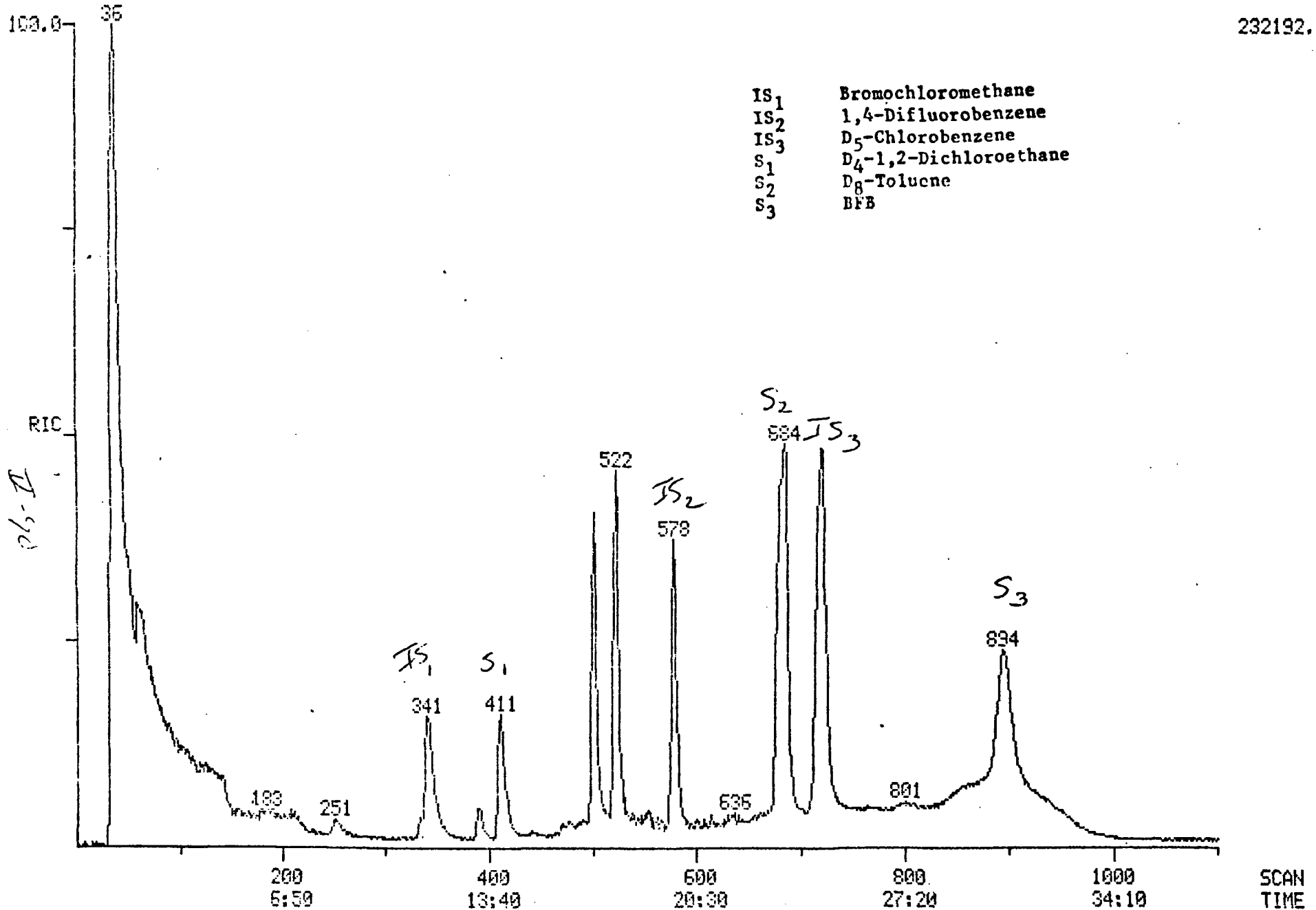
RIC
03/05/85 14:43:00
SAMPLE: 43232 MSD CASE 3963 AB-028 MSD

DATA: U2742

SCANS 1 TO 1100

232192.

IS ₁	Bromochloromethane
IS ₂	1,4-Difluorobenzene
IS ₃	D ₅ -Chlorobenzene
S ₁	D ₄ -1,2-Dichloroethane
S ₂	D ₈ -Toluene
S ₃	BFB



DATA: V2742.TI
 03/05/85 14:43:00
 SAMPLE: 43232 MSD CASE 3569 AB-028 MSD
 SUBMITTED BY: ANALYST:

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

- NO NAME
- 1 BROMOCHLOROMETHANE (I. S. #1)
- 2 CHLOROMETHANE
- 3 BROMOMETHANE
- 4 VINYL CHLORIDE
- 5 CHLOROETHANE
- 6 METHYLENE CHLORIDE
- 7 ACETONE
- 8 CARBON DISULFIDE
- 9 1, 1-DICHLOROETHENE
- 10 1, 1-DICHLOROETHANE
- 11 TRANS 1, 2-DICHLOROETHENE
- 12 CHLOROFORM
- 13 1, 2-DICHLOROETHANE
- 14 D4-1, 2-DICHLOROETHANE
- 15 1, 1, 1-TRICHLOROETHANE
- 16 1, 4-DIFLUOROBENZENE (I. S. #2)
- 17 CARBON TETRACHLORIDE
- 18 VINYL ACETATE
- 19 2-BUTANONE
- 20 BROMODICHLOROMETHANE
- 21 1, 2-DICHLOROPROPANE
- 22 TRANS 1, 3-DICHLOROPROPANE
- 23 TRICHLOROETHENE
- 24 DIBROMOCHLOROMETHANE
- 25 1, 1, 2-TRICHLOROETHANE
- 26 BENZENE
- 27 CIS-1, 3-DICHLOROPROPENE
- 28 BROMOFORM
- 29 D8-TOLUENE (SURRE#2)
- 30 2-CHLOROETHYL VINYLETHER
- 31 D5-CHLOROBENZENE (I. S. #3)
- 32 4 METHYL 2 PENTANONE
- 33 2-HEXANONE
- 34 TETRACHLOROETHENE
- 35 1, 1, 2, 2, TETRACHLOROETHANE
- 36 TOLUENE
- 37 CHLOROBENZENE
- 38 ETHYLBENZENE
- 39 BFF (SURRE#3)
- 40 STYRENE
- 41 TOTAL XYLENES

NO	M/E	SCAN	TIME	REF	RRT	METH	AREA(HGHT)	AMOUNT	%TOT
1	128	341	11:39	1	1.000	A BB	29442.	50.000 UG/L	7.60
2	NOT FOUND								
3	NOT FOUND								

11-91

NO	SIZE	SCAN	TIME	REF	VAL	UNIT	OPERATOR	AMOUNT	UNIT	WGT
4		NOT FOUND								
5		NOT FOUND								
6		NOT FOUND								
7		NOT FOUND								
8		NOT FOUND								
9		NOT FOUND								
10		NOT FOUND								
11		NOT FOUND								
12	83	390	13:19	1	1.144	A BB	22286.	7.163	UG/L	1.09
13		NOT FOUND								
14	65	411	14:03	1	1.205	A BB	99907.	93.593	%REC	14.22
15		NOT FOUND								
16	114	578	19:45	16	1.000	A BB	148270.	50.000	UG/L	7.60
17		NOT FOUND								
18		NOT FOUND								
19		NOT FOUND								
20		NOT FOUND								
21		NOT FOUND								
22		NOT FOUND								
23	130	501	17:07	16	0.867	A BB	63763.	55.794	UG/L	8.48
24		NOT FOUND								
25		NOT FOUND								
26	78	522	17:50	16	0.903	A BB	215462.	53.146	UG/L	8.07
27		NOT FOUND								
28		NOT FOUND								
29	98	679	23:12	16	1.175	A BB	174924.	97.031	%REC	14.74
30		NOT FOUND								
31	117	716	24:28	31	1.000	A BB	109539.	50.000	UG/L	7.60
32		NOT FOUND								
33		NOT FOUND								
34	164	634	21:40	31	0.885	A BB	577.	0.585	UG/L	0.09
35		NOT FOUND								
36	92	685	23:24	31	0.957	A BB	133973.	53.661	UG/L	8.15
37	112	720	24:36	31	1.006	A BB	139944.	56.295	UG/L	8.55
38		NOT FOUND								
39	95	894	33:33	31	1.249	A BB	88957.	90.921	%REC	13.81
40		NOT FOUND								
41		NOT FOUND								

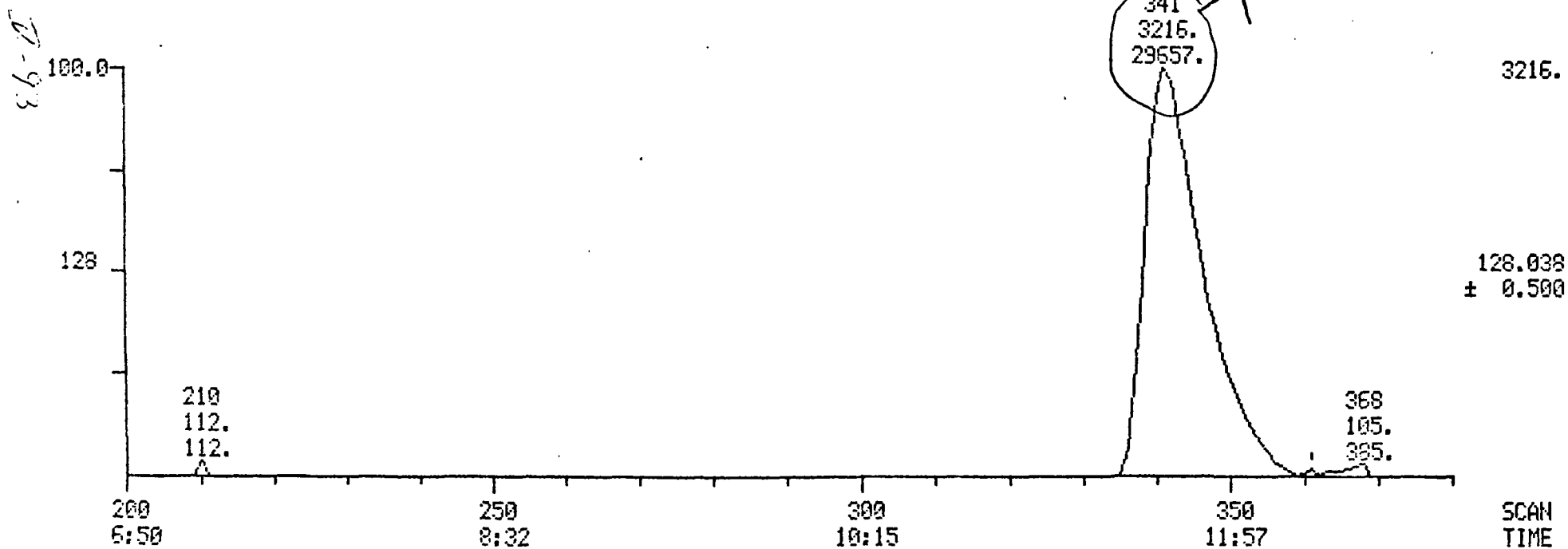
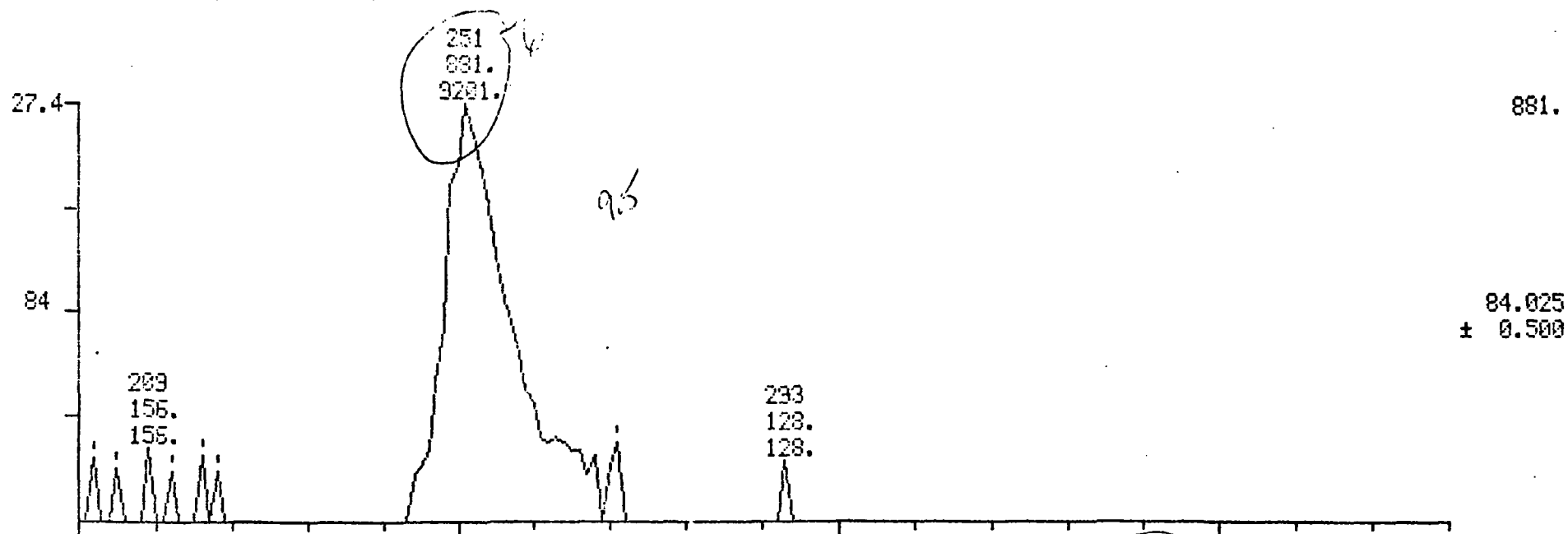
AB 028 MSD Case 3969 3/5/85

D-92

MASS CHROMATOGRAMS
03/05/85 14:43:00
SAMPLE: 43232 MSD CASE 3969 AB-028 MSD

DATA: U2742

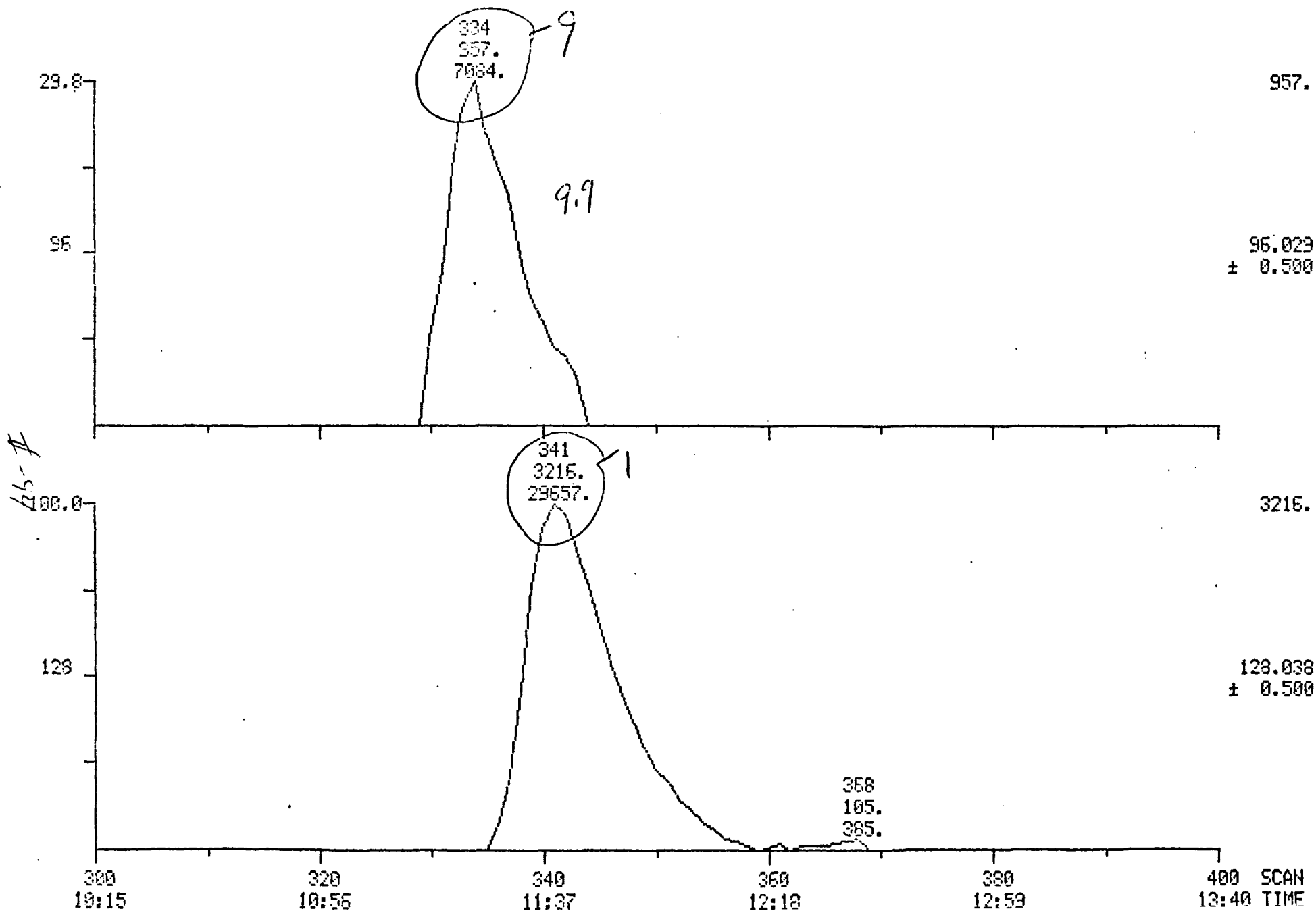
SCANS 200 TO 380



MASS CHROMATOGRAMS
03/05/85 14:43:00
SAMPLE: 43232 MSD CASE 3959 AB-028 MSD

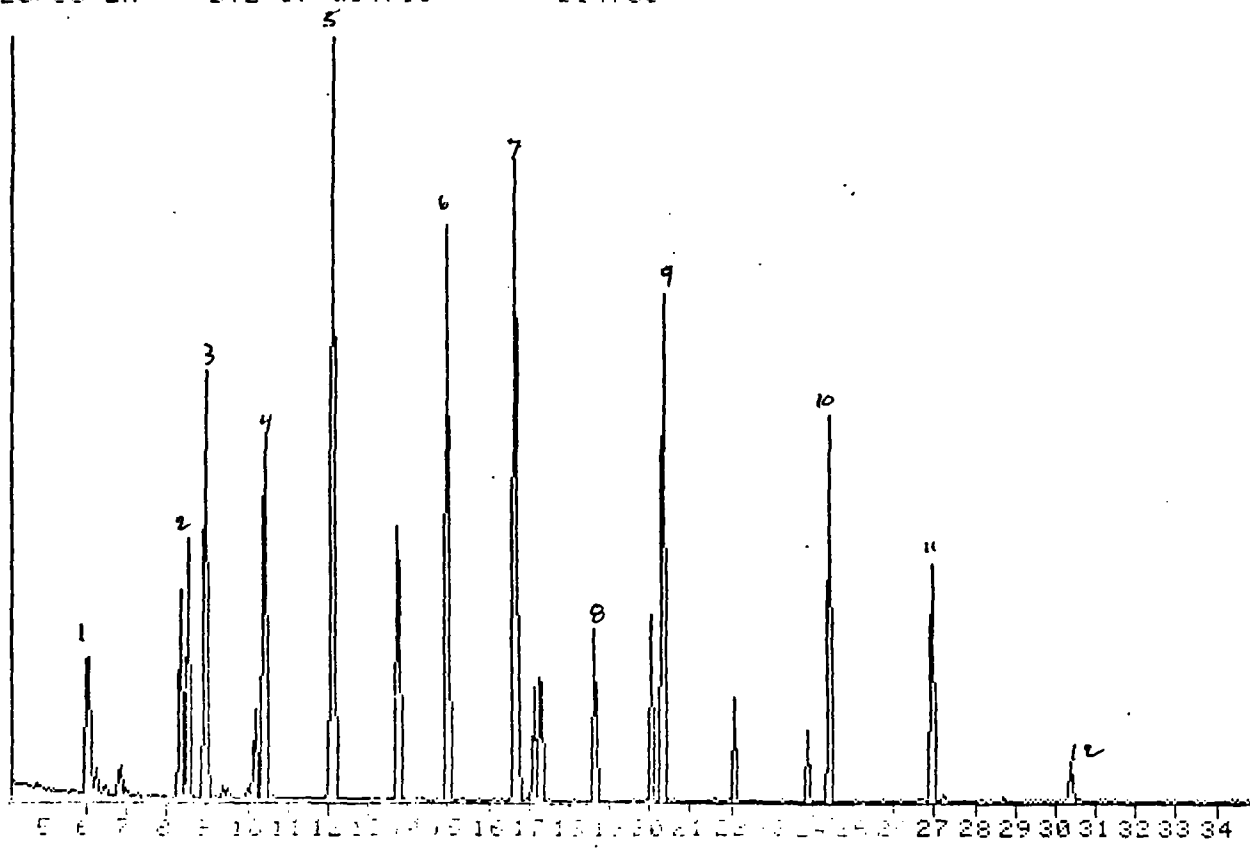
DATA: V2742

SCANS 300 TO 400



81828

4-10



1. 2-FLUOROPHENOL (SUFR)
2. PHENOL-D6 (SUFR)
3. 1,4-DICHLOROBENZENE-D4 (IS)
4. NITROBENZENE-D5 (SUFR)
5. NAPHTHALENE-D8 (IS)
6. 2-FLUORODIPHENYL (SUFR)
7. ACENAPHTHENE-D10 (IS)
8. 2,4,6-TRIBROMOPHENOL (SUFR)
9. PHENANTHRENE-D10 (IS)
10. TERPHENYL-D14 (SUFR)
11. CHRYSENE-D12 (IS)
12. PERYLENE-D12 (IS)

D-95

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
9.0	150.0	1.000	80.0000				D4-1,4-DICHLOROBENZENE(20)
9.9	74.0	.046	.0000				N-NITROSODIMETHYLAMINE
8.4	93.0	.913	.0000				ANILINE
8.5	93.0	.709	1.7247 ^{EM}				BIS(2-CHLOROETHYL)ETHER
8.9	146.0	.591	.0000				1,3-DICHLOROBENZENE
9.0	146.0	.594	42.6409 [✓]				1,4-DICHLOROBENZENE
9.5	108.0	.395	.0000				BENZYL ALCOHOL
9.5	146.0	.594	.0000				1,2-DICHLOROBENZENE
9.9	45.0	.527	.0000				BIS(2-CHLOROISOPROPYL)ETHER
10.2	70.0	.490	54.0577 [✓]				N-NITROSODINPROPYLAMINE
10.2	117.0	.291	.0000				HEXACHLOROETHANE
10.5	77.0	.716	.0000				NITROBENZENE

12.1	68.0	1.000	80.0000				D8-NAPHTHALENE(20)
11.1	138.0	1.519	.0000				ISOPHORONE
11.8	95.0	1.813	.0000				BIS(2-CHLOROETHOXY)METHANE
12.1	180.0	2.642	39.6608 [✓]				1,2,4-TRICHLOROBENZENE
12.2	129.0	1.325	.0000				NAPHTHALENE
12.5	127.0	4.577	.0000				4-CHLOROANILINE
12.7	225.0	1.190	.0000				HEXACHLOROBTADIENE
14.0	115.0	2.578	.0000				2-METHYLNAPHTHALENE

16.7	164.0	1.000	80.0000				D10-ACENAPHTHENE
14.6	237.0	.205	.0000				HEXACHLOROCYCLOPENTADIENE
15.2	164.0	.281	.0000				2-CHLORONAPHTHALENE
15.6	138.0	.397	.0000				2-NITROANILINE
16.2	164.0	.083	.0000				DIMETHYL PHTHALATE
16.3	151.0	.272	.0000				ACENAPHTHYLENE
16.4	165.0	.228	.0000				2,6-DINITROTOLUENE
16.7	138.0	.293	.0000				3-NITROANILINE
16.7	152.0	.437	43.2235 [✓]				ACENAPHTHENE
17.2	139.0	.490	24.2522 ^{EM}				DIBENZOFURAN
17.3	165.0	.320	44.0576 [✓]				2,4-DINITROTOLUENE
18.0	177.0	.171	.0000				DIETHYL PHTHALATE
18.0	165.0	.820	.0000				FLUORENE
18.1	204.0	.358	.0000				4-CHLOROPHENYLPHENYL ETHER
18.2	138.0	.292	.0000				4-NITROANILINE
17.2	169.0	.144	.0000				DIPHENYLAMINE

20.3	188.0	1.000	80.0000				D10-PHENANTHRENE(20)
19.3	248.0	.112	.0000				4-BROMOPHENYLPHENYL ETHER
19.6	284.0	.112	.0000				HEXACHLOROBENZENE
20.4	179.0	.060	.0000				PHENANTHRENE/ANTHRACENE
22.1	150.0	.100	16.6396 [✓]				DIBUTYL PHTHALATE
23.4	202.0	.665	.0000				FLUORANTHENE

D-96

FILE NUMBER 14758
 QC-958 AB028MSD CASE 3969
 3/25/85/EM BTL#17 Q14758 D14758

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
27.0	240.0	1.000	80.0000				D12-CHRYSENE
23.9	202.0	2.264	12.6747	✓			PYRENE
25.8	149.0	1.136	.0000				BUTYL BENZYL PHTHALATE
26.9	228.0	.427	.0000				BENZO(A)ANTHRACENE/CHRYSENE
27.2	149.0	1.128	5.1470	EM			BIS(2-ETHYLHEXYL)PHTHALATE

30.4	264.0	1.000	80.0000				D12-PERYLENE
28.7	149.0	5.974	2.5719	EM			DIOCTYL PHTHALATE
29.5	252.0	1.307	.0000				BENZO(B)/(K)FLUORANTHENE
30.2	252.0	.744	.0000				BENZO(A)PYRENE
33.7	276.0	.485	.0000				INDENO(123-CD)PYRENE
33.8	278.0	.346	.0000				DIBENZO(AH)ANTHRACENE
34.6	276.0	.360	.0000				BENZO(GHI)PERYLENE

27.0	240.0	1.000	80.0000				D12-CHRYSENE
23.8	184.0	.712	.0000				BENZIDINE
27.0	252.0	.146	.0000				DICHLOROBENZIDINE

D-97

AREA	R. TIME	START	STOP	MAXIMA	# SUMS	MASS	BASE
0	150.0	1.000	80.0000	✓	D4-1,4-DICHLOROBENZENE(20)		
4	94.0	.841	51.8325	✓	PHENOL		
5	128.0	.611	85.0818	✓	2-CHLOROPHENOL		
9	108.0	.671	.0000		2-METHYLPHENOL		
2	108.0	.699	.0000		4-METHYLPHENOL		

1	68.0	1.000	80.0000		D8-NAPHTHALENE(20)		
3	139.0	1.759	.0000		2-NITROPHENOL		
5	122.0	2.581	.0000		2,4-DIMETHYLPHENOL		
1	122.0	1.200	.0000		BENZOIC ACID		
9	162.0	2.051	.0000		2,4-DICHLOROPHENOL		
7	142.0	2.196	83.5878	✓	4-CHLORO-M-CRESOL		

7	164.0	1.000	80.0000		D10-ACENAPHTHENE(20)		
8	196.0	.237	.0000		2,4,5-TRICHLOROPHENOL		
8	196.0	.237	.0000		2,4,6-TRICHLOROPHENOL		
9	184.0	.111	.0000		2,4-DINITROPHENOL		
3	198.0	.126	.0000		4,6-DINITRO-O-CRESOL		
2	65.0	.231	53.8130	✓	4-NITROPHENOL		

3	188.0	1.000	80.0000		D10-PHENANTHRENE		
1	266.0	.072	76.2480	✓	PENTACHLOROPHENOL		

Date 4/1/75

Analyst BL

Contract 3-705-050

Pesticide Analysis - Qualitative/Quantitative

Case # 3769

Sample I.D. CC-968/ABO30 MSD
Run # 16

HP 58 80
Bottle # 16

RT (min.)	Peak ID	Area x 10 ³	ng/ml	Dilution Factor (mls extracted)	Spiked at ng	% Recovery (ug/l) ppb	total ug
2.57	δ -BHC	29.49	*13.9 x 10	=139	200	69.5	
3.18	Heptachlor	35.58	13.6 x 10	=136	200	68.0	
3.87	Aldrin	32.35	15.9 x 10	=159	200	79.5	
9.42	Dieldrin	65.65	39.9 x 10	=399	500	79.8	
11.52	Endrin	67.12	56.8 x 10	=568	500	114	
17.37	pp'-DDT	42.71	56.6 x 10	=566	500	113	
	DBC obscured by peak at 33 min						
	* Calculated from lowest four calibration points						
	* Calculated from highest four calibration points						

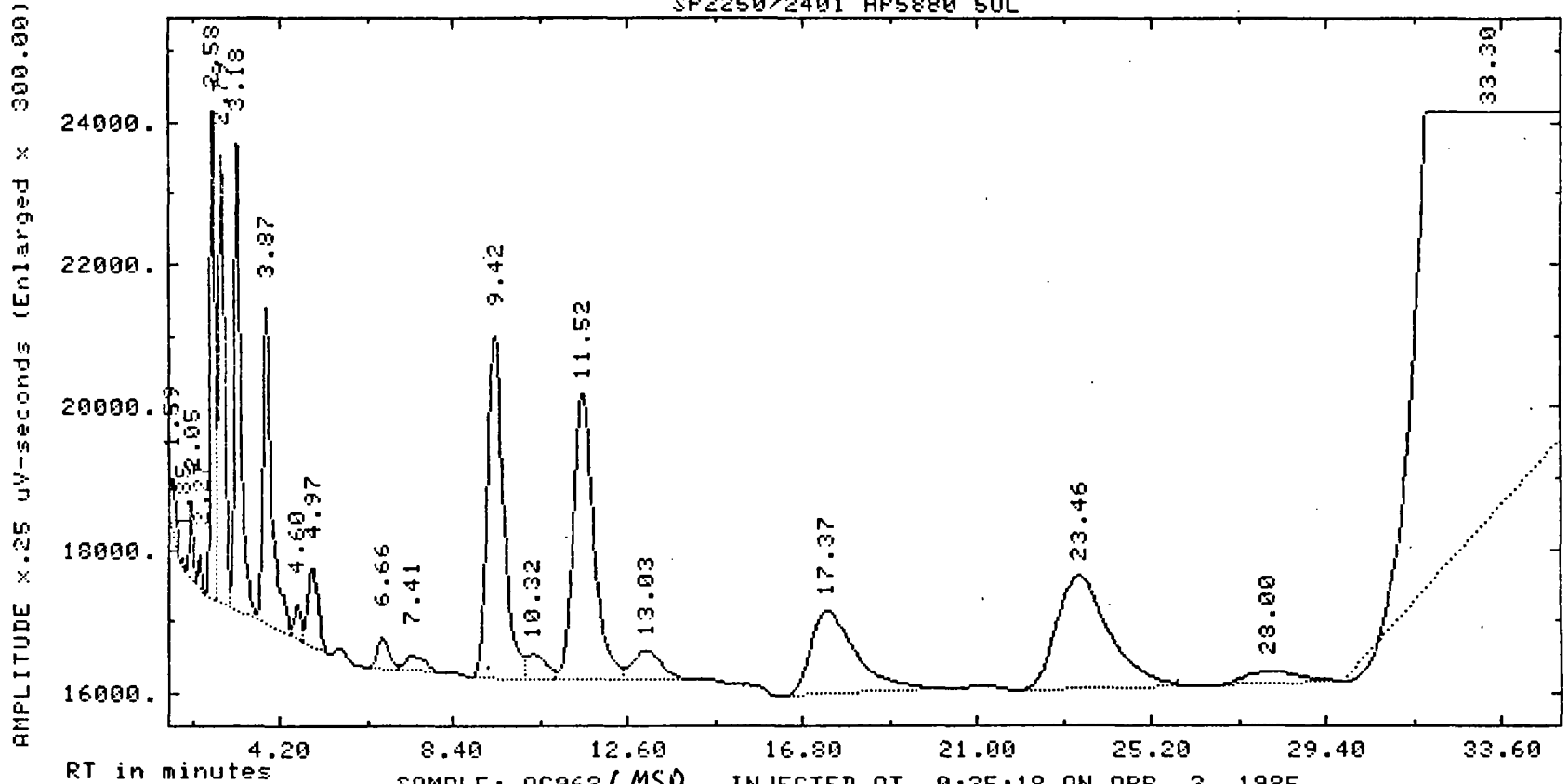


GCA CORPORATION
Technology Division

II-99

Case 3969

SP2250/2401 HP5880 5UL



SAMPLE: QC968/MSD INJECTED AT 0:25:18 ON APR 2, 1985
Method: PRIME Raw: RB5016 Proc: PB5016

001-11

APR 2, 1985 9:14:15

(USE 010)

REPORT: 159 CHANNEL: 2 # PEAKS: -24 SP2250/2401 HP5080 SUL

PEAK	RT	ITH	FACTOR	AREA	AREA %	NAME
1	1.10	0.00	1.00000	4104 BV	.123	
2	1.16	0.00	1.00000	5421 VV	.163	
3	1.37	0.00	1.00000	5070 VV	.153	
4	1.59	0.00	1.00000	2765 VV	.084	
5	1.85	0.00	1.00000	250 VV	.008	
6	2.05	0.00	1.00000	3605 VV	.108	
7	2.27	0.00	1.00000	1425 VV	.043	
8	2.50	0.00	1.00000	29487 VV	.887	<i>δ-BHC</i>
9	2.79	0.00	1.00000	30028 VV	.904	
10	3.18	0.00	1.00000	35570 VV	1.071	<i>Heptachlor</i>
11	3.87	0.00	1.00000	32349 VV	.973	<i>Aldrin</i>
12	4.60	0.00	1.00000	2742 VV	.083	
13	4.97	0.00	1.00000	9131 VD	.275	
14	6.66	0.00	1.00000	4865 BV	.146	
15	7.41	0.00	1.00000	3321 VD	.100	
16	9.42	0.00	1.00000	65852 BV	1.976	<i>Dieldrin</i>
17	10.32	0.00	1.00000	5462 VV	.164	
18	11.52	0.00	1.00000	67117 VV	2.020	<i>Endrin</i>
19	13.03	0.00	1.00000	9153 VB	.276	
20	17.37	0.00	1.00000	42814 BB	1.288	<i>pp'-DDT</i>
21	23.46	0.00	1.00000	69270 BV	2.085	
22	26.03	0.00	1.00000	8140 VD	.245	
23	33.30	0.00	1.00000	2857193 BV	85.979	
24	39.45	0.00	1.00000	28132 VF	.847	

TOTAL AREA = 3323114 TOTAL AREA % = 100.000

SAMPLE: 00906/MSD *AR30* INJECTED AT 8:25:18 ON APR 2, 1985

ZERO METHOD: PRIME SEQ: SMO3A SUBSQ/SAMP: 1 / 16 BTL: 10

SL-UBTH .500 MV/MIN .300 DELAY 1.00 MIN-AR 50 BUNCH AUTO REPORT MEDIUM

SUP-UNK NO DVT 0.00 ID-LVL 0 REF-RTW .30 %RTW 5.0 %DIL-F 100.00 ISO YES

ACTUAL RUN TIME: 40.000 MINUTES

RF-UNK: 1.00000 STD-AMT: 1.0000 SAMP-AMT: 1.0000

ENDED NOT ON BASELINE

RAW DATA FILE: RBS016 PARAM FILES= METHOD: SEQ:

DONE

II-101

CAMP DRESSER & MCKEE INC.

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SUITE 500

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703 642-5500