

*Chlorophenols.*

SAMPLE DATA PACKET

010338

U.S. ENVIRONMENTAL PROTECTION AGENCY  
 CLP Sample Management Office  
 P.O. Box 818 - Alexandria, Virginia 22313  
 Phone: 703/557-2490 - FTS/557-2490

SAS Number  
**2493-F**

**SPECIAL ANALYTICAL SERVICE  
 PACKING LIST**

<b>Sampling Office:</b> <u>Region 6</u>	<b>Sampling Date(s):</b> <u>10-2-86</u>	<b>Ship To:</b> <u>ENVIRODYNE ENGINEERS</u> <u>12161 LACKLAND RD</u> <u>ST LOUIS, MO 63146</u>	<b>For Lab Use Only</b>
<b>Sampling Contact:</b> <u>Dave Anderson</u> (name)	<b>Date Shipped:</b> <u>10-2-86</u>	<b>Attn:</b> <u>DR. MARGARET C. WINTER</u>	<b>Date Samples Rec'd:</b> <u>10-3-86</u>
<u>(214) 742-6601</u> (phone)	<b>Site Name/Code:</b> <u>[REDACTED]</u>		<b>Received By:</b> <u>[Signature]</u>

Sample Numbers	Sample Description i.e., Analysis, Matrix, Concentration	Sample Condition on Receipt at Lab
1. <u>2493F01</u>	<u>Low dust wipe TCDD Sta 01</u>	<u>GOOD</u>
2. <u>2493F02</u>	<u>Low dust wipe TCDD Sta 02</u>	
3. <u>2493F03</u>	<u>Low dust wipe TCDD Sta 03</u>	
4. <u>2493F04</u>	<u>Low dust wipe TCDD Sta 04</u>	
5. <u>2493F05</u>	<u>Low dust wipe TCDD MSDD</u>	
6. <u>2493F06</u>	<u>Low dust wipe TCDD MSDD</u>	
7. <u>2493F07</u>	<u>Low dust wipe TCDD MSDD</u>	
8. <u>2493F08</u>	<u>Low dust wipe chlorinated phenols Sta 01</u>	
9. <u>2493F09</u>	<u>Low dust wipe chlorinated phenols Sta 02</u>	
10. <u>2493F10</u>	<u>Low dust wipe chlorinated phenols Sta 03</u>	
11. <u>2493F11</u>	<u>Low dust wipe chlorinated phenols Sta 04</u>	
12. <u>2493F12</u>	<u>Low dust wipe chlorinated phenols MSAP</u>	
13. <u>2493F13</u>	<u>Low dust wipe chlorinated phenols MSAP</u>	
14. <u>2493F14</u>	<u>Low dust wipe chlorinated phenols MSOP</u>	
15.		
16.		
17.		
18.		
19.		
20.		

0103010

**For Lab Use Only**

White - SMO Copy, Yellow - Region Copy, Pink - Lab Copy for return to SMO, Gold - Lab Copy

Organics Analysis Data Sheet  
(Page 1)

Laboratory Name: ENVIRONMENTAL ENGRS Case No: 3101-2493F  
 Lab Sample ID No: 86002739 QC Report No: \_\_\_\_\_  
 Sample Matrix: WIPES Contract No: SAS 3101-2493F  
 Date Release Authorized By: MARCOLET C. WILSON Date Sample Received: 10-6-86

## Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: \_\_\_\_\_

Date Analyzed: \_\_\_\_\_

Conc/Dil Factor: \_\_\_\_\_ pH \_\_\_\_\_

Percent Moisture: (Not Decanted) \_\_\_\_\_

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	
67-64-1	Acetone	
75-15-0	Carbon Disulfide	
75-35-4	1, 1-Dichloroethane	
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	

CAS Number		ug/l or ug/Kg (Circle One)
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	
75-35-2	Bromoforn	
108-10-1	4-Methyl-2-Pentanone	
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	
108-90-7	Chlorobenzene	
100-41-4	Ethylbenzene	
100-42-5	Styrene	
	Total Xylenes	

## Data Reporting Qualifiers

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

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10J). If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3J.
- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/ul in the final extract should be confirmed by GC/MS.
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Laboratory Name: ENVIRODYNE

Case No: 2493F

Sample Number  
2493F-08

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration: (Low) Medium (Circle One)

Date Extracted/Prepared: 10/6/86

Date Analyzed: 10/16/86

Conc/Dil Factor: 10

Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No

Separatory Funnel Extraction  Yes

Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/wipe ug/100ug/Kg (Circle One)
108-95-2	Phenol	
111-44-4	bis(-2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	1000
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	
120-83-2	2, 4-Dichlorophenol	1000
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	1000
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2, 4, 6-Trichlorophenol	1000
95-95-4	2, 4, 5-Trichlorophenol	5000
91-58-7	2-Chloronaphthalene	
88-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
206-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	

CAS Number		ug/wipe ug/100ug/Kg (Circle One)
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	
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	5000
85-01-8	Phenanthrene	
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
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	
218-01-9	Chrysene	
117-84-0	Di-n-Octyl Phthalate	
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	Dibenz(a, h)Anthracene	
191-24-2	Benzo(g, h, i)Perylene	

(1)-Cannot be separated from diphenylamine

108010

Laboratory Name: ENVIRODYNE

Case No: 2493F

Sample Number

2493F-08

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.	no chlorophenols found			
2.				
3.				
4.				
5.				
6.				
7.				
8.				
9.				
10.				
11.				
12.				
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14.				
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21.				
22.				
23.				
24.				
25.				
26.				
27.				
28.				
29.				
30.				

010300

QUANT REPORT

Operator ID: USER8      Quant Rev: 4      Quant Time: 861020 12:34  
 Output File: ^A5257::L2      Injected at: 861016 05:02  
 Data File: >A5257::E2      Dilution Factor: 1.000 10      mcv 10-21-86  
 Name: 2493F-08  
 Misc: LMC 5996 86002739 3101-2493F      BTL#24

ID File: #A5248::03  
 Title: ID FILE FOR SAS 3101-2493F  
 Last Calibration: 861020 11:54

Compound	R.T.	Scan#	Area	Conc	Units	q
1) #D4 1,4-DICHLOROBENZENE	8.18	490	143960	40.00	UG/WIPE	94
2) 2-FLUOROPHENOL (SURROGATE)	5.33	213	20634	<del>5.91</del>	UG/WIPE	100
4) #D8 NAPHTHALENE	11.83	845	347300	40.00	UG/WIPE	100
7) #D10 ACENAPHTHENE	17.25	1370	183112	40.00	UG/WIPE	100
8) 2,4,6-TRIBROMOPHENOL (SURR)	19.72	1609	14202	<del>16.74</del>	UG/WIPE	100
11) #D10 PHENANTHRENE	21.78	1800	459705	40.00	UG/WIPE	100

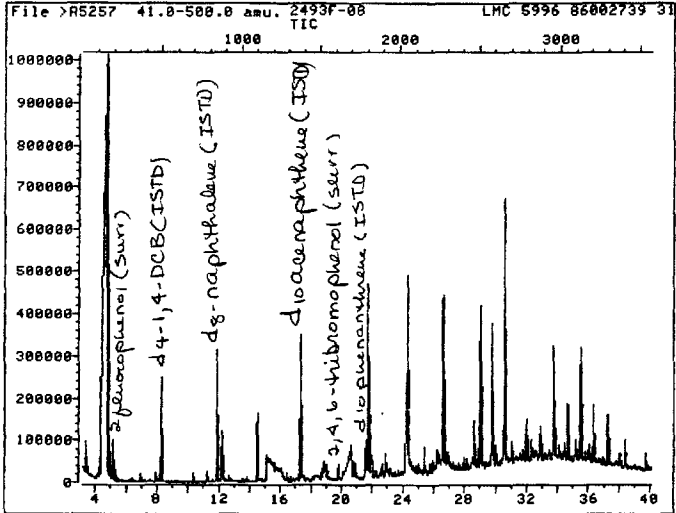
99.1 mcv 10-21-86

167.4 mcv 10-21-86

\* Compound is ISTD

0108910

TOTAL ION CHROMATOGRAM



Data File: >A5257::E2  
Name: 2493F-08  
Misc: LMC 5996 86002739 3101-2493F

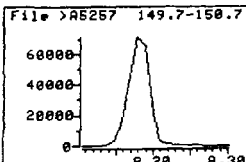
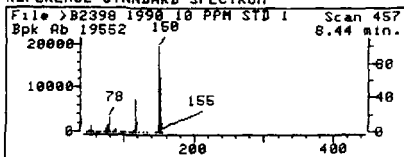
BTL424

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Last Calibration: 861020 11:54

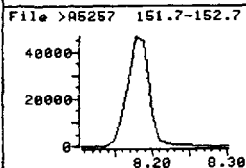
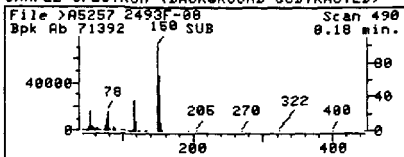
Operator ID: USERS  
Quant Time: 861020 12:34  
Injected at: 861016 05:02

010234

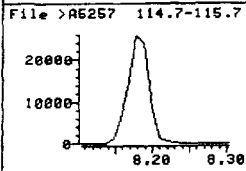
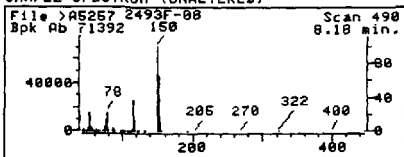
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



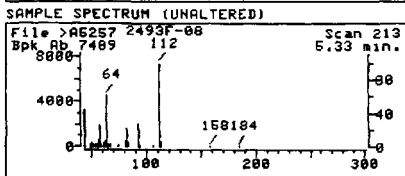
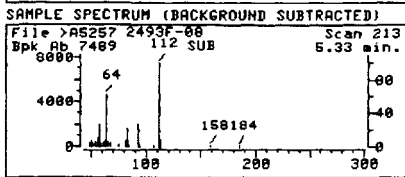
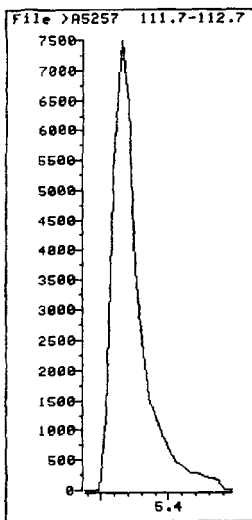
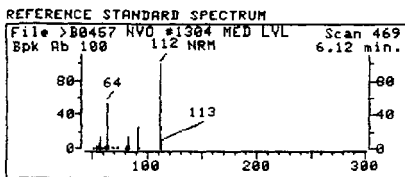
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 Name: 2493F-08  
 Misc: LMC 5996 86002739 3101-2493F  
 Quant Time: 861020 12:34  
 Injected at: 861016 05:02

BTL424

Compound No: 1 (ISTD)  
 Compound Name: D4 1,4-DICHLOROBENZENE  
 Scan Number: 490  
 Retention Time: 8.18 min.  
 Area: 143960  
 Concentration: 40.00 UG/WIPE  
 q-value: 94

010305





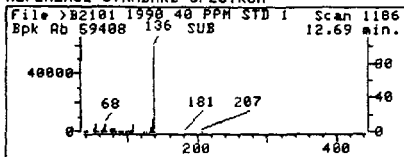
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Injected at: 861016 05:02

BTL#24

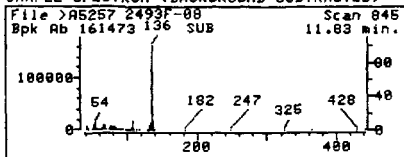
Compound No: 2  
Compound Name: 2-FLUOROPHENOL (SURROGATE)  
Scan Number: 213  
Retention Time: 5.33 min.  
Area: 20634  
Concentration: 9.51 UG/WIPE 99.1 mcw 10/21/86  
q-value: 100

010326

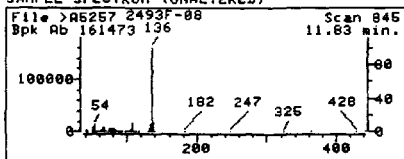
REFERENCE STANDARD SPECTRUM



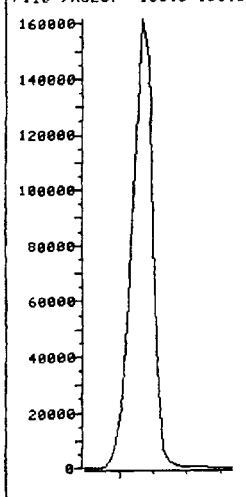
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SAMPLE SPECTRUM (UNALTERED)



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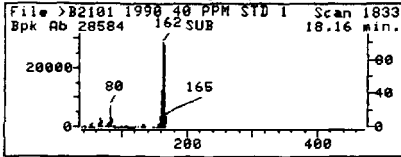
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Injected at: 861016 05:02

BTL424

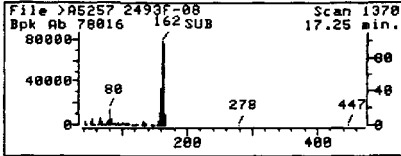
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Compound Name: 08 NAPHTHALENE  
Scan Number: 845  
Retention Time: 11.83 min.  
Area: 347300  
Concentration: 40.00 UG/WIPE  
q-value: 100

010307

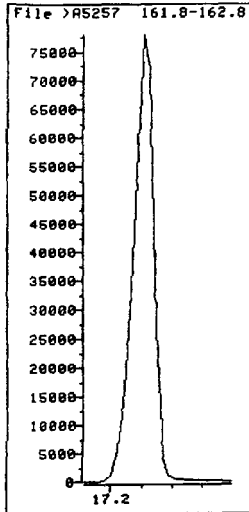
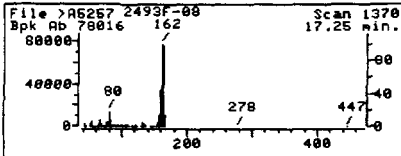
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

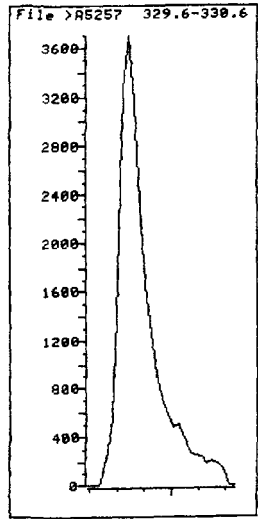
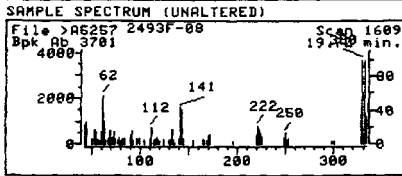
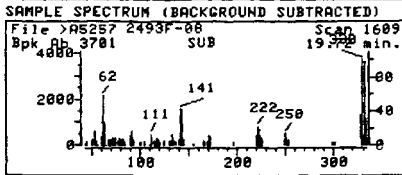
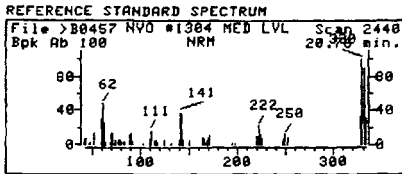


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Name: 2493F-08  
Misc: LMC 5996 06002739 3101-2493F  
Quant Time: 061020 12:34  
Injected at: 061016 05:02

BT424

Compound No: 7 (ISTD)  
Compound Name: D10 ACENAPHTHENE  
Scan Number: 1370  
Retention Time: 17.25 min.  
Area: 183112  
Concentration: 40.00 UG/WIPE  
q-value: 100

010308



Data File: >A5257::E2  
Name: 2493F-08  
Misc: LMC 5996 86002739 3101-2493F  
Quant Time: 861020 12:34  
Injected at: 861016 05:02

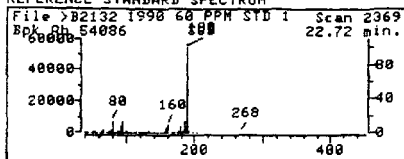
BTL#24

Compound No: 8  
Compound Name: 2,4,6-TRIBROMOPHENOL (SURR)  
Scan Number: 1609  
Retention Time: 19.72 min.  
Area: 14202  
Concentration: 16.74 ~~UG/WIPE~~  
q-value: 100

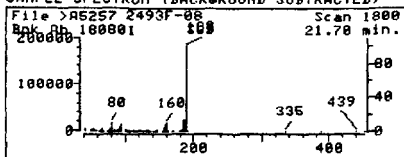
167.4  $\mu\text{g/wipe}$  mow 10-21-86

010300

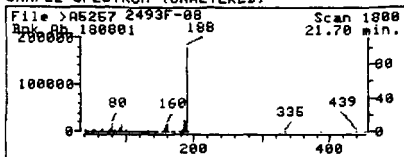
REFERENCE STANDARD SPECTRUM



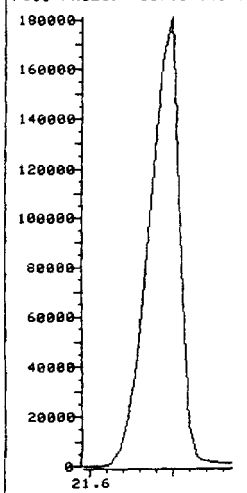
SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



File >A5257 187.8-188.8



Data File: >A5257::E2  
Name: 2493F-08  
Misc: LMC 5996 86002739 3101-2493F  
Quant Time: 861020 12:34  
Injected at: 861016 05:02

BTL424

Compound No: 11 (ISTD)  
Compound Name: D10 PHENANTHRENE  
Scan Number: 1800  
Retention Time: 21.70 min.  
Area: 459705  
Concentration: 40.00 UG/WIPE  
q-value: 100

010370

## Organics Analysis Data Sheet

(Page 1)

Laboratory Name: ENVIRONMENTAL ENGRS. Case No: 3101-2493F  
 Lab Sample ID No: 86002740 QC Report No: \_\_\_\_\_  
 Sample Matrix: WIPE Contract No: SAS 3101-2493F  
 Data Release Authorized By: Margaret C. Wille Date Sample Received: 10/6/86

## Volatile Compounds

Concentration: (Low) Medium (Circle One)

Date Extracted/Prepared: \_\_\_\_\_

Date Analyzed: \_\_\_\_\_

Conc/Dil Factor: \_\_\_\_\_ pH \_\_\_\_\_

Percent Moisture: (Not Decanted) \_\_\_\_\_

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	
67-64-1	Acetone	
75-15-0	Carbon Disulfide	
75-35-4	1, 1-Dichloroethane	
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	

CAS Number		ug/l or ug/Kg (Circle One)
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	
108-10-1	4-Methyl-2-Pentanone	
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	
108-90-7	Chlorobenzene	
100-41-4	Ethylbenzene	
100-42-5	Styrene	
	Total Xylenes	

## Data Reporting Qualifiers

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

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10J). If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 3J.
- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides 2:10 ug/l of in the final extract should be confirmed by GC/MS.
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Laboratory Name: ENRODYNE

Case No: 2493F

Sample Number  
2493F-09

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration: (Low) Medium (Circle One)  
Date Extracted/Prepared: 10/6/86  
Date Analyzed: 10-16-86  
Conc/Dil Factor: 1  
Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No  
Separatory Funnel Extraction  Yes  
Continuous Liquid - Liquid Extraction  Yes

CAS Number	Compound	ug/wipe ug/l or ug/kg (Circle One)
108-95-2	Phenol	
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	100
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	
120-83-2	2, 4-Dichlorophenol	100
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	100
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2, 4, 6-Trichlorophenol	100
95-95-4	2, 4, 5-Trichlorophenol	500
91-58-7	2-Chloronaphthalene	
88-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	

CAS Number	Compound	ug/wipe ug/l or ug/kg (Circle One)
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	
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	500
85-01-8	Phenanthrene	
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
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	
218-01-9	Chrysene	
117-84-0	Di-n-Octyl Phthalate	
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	Dibenz(a,h)Anthracene	
191-24-2	Benzo(g,h,i)Perylene	

(1)-Cannot be separated from diphenylamine

010370

Laboratory Name: ENV RODYNE  
Case No: 2493F

Sample Number  
2493F-09

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 chloro phenols found</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.				

010379



QUANT REPORT

Operator ID: USER8                      Quant Rev: 4    Quant Time: 861020 12:30  
 Output File: ^A5254::L2                      Injected at: 861016 02:40  
 Data File: >A5254::E2                      Dilution Factor: 1.000  
 Name: 2493F-09  
 Misc: LMC 5996 86002740 3101-2493F                      BTL#21

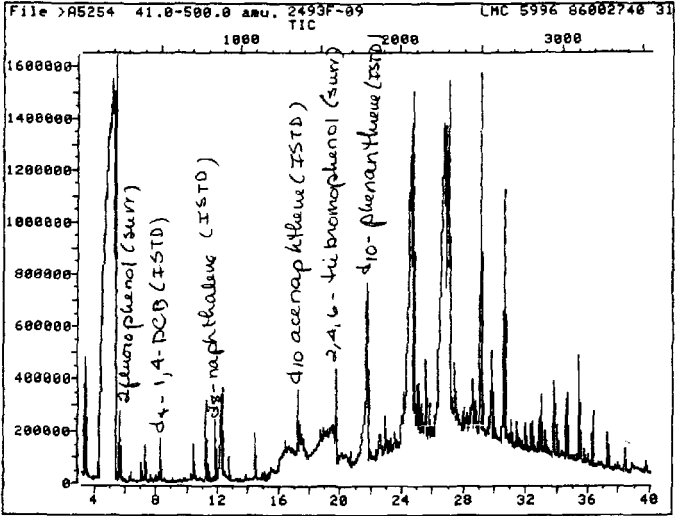
ID File: #A5248::03  
 Title: ID FILE FOR SAS 3101-2493F  
 Last Calibration: 861020 11:54

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *D4 1,4-DICHLOROBENZENE	8.17	491	109098	40.00	UG/WIPE	95
2) 2-FLUOROPHENOL (SURROGATE)	5.59	241	135952	86.13	UG/WIPE	100
4) *D8 NAPHTHALENE	11.81	844	246434	40.00	UG/WIPE	100
7) *D10 ACENAPHTHENE	17.20	1365	129711	40.00	UG/WIPE	100
8) 2,4,6-TRIBROMOPHENOL (SURR)	19.70	1606	119123	198.19	UG/WIPE	100
11) *D10 PHENANTHRENE	21.65	1793	389674	40.00	UG/WIPE	100

\* Compound is ISTD

718070

TOTAL ION CHROMATOGRAM



Data File: >A5254::E2  
Name: 2493F-09  
Misc: LMC 5996 86002740 3101-2493F

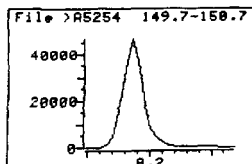
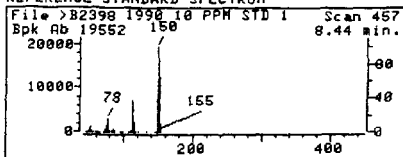
BT1#21

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Title: ID FILE FOR SAS 3101-2493F  
Last Calibration: 861020 11:54

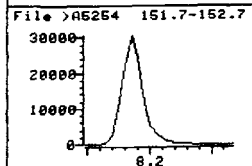
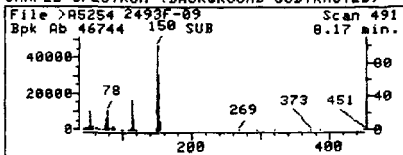
Operator ID: USER8  
Quant Time: 861020 12:30  
Injected at: 861016 02:40

010375

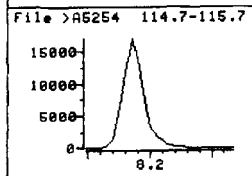
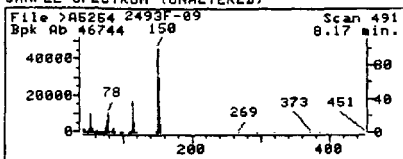
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



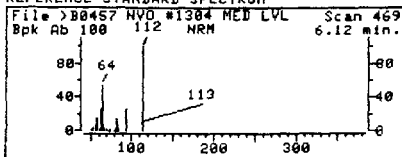
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Name: 2493F-09  
Misc: LMC 5996 86002740 3101-2493F  
Quant Time: 861028 12:30  
Injected at: 861016 02:40

BTL#21

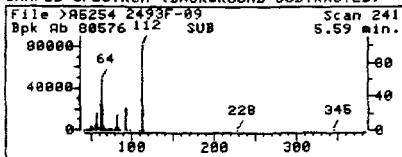
Compound No: 1 (ISTD)  
Compound Name: D4 1,4-DICHLOROBENZENE  
Scan Number: 491  
Retention Time: 8.17 min.  
Area: 109098  
Concentration: 40.00 UG/WIPE  
q-value: 95

0103070

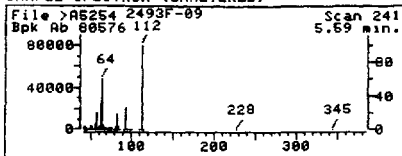
REFERENCE STANDARD SPECTRUM



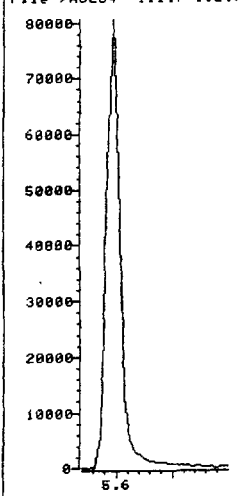
SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



File >A6254 111.7-112.7



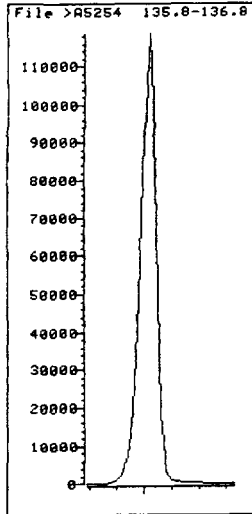
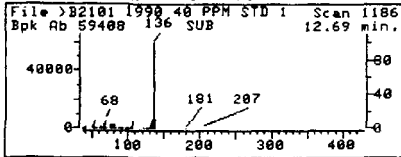
Data File: >A6254::E2  
Name: 2493F-09  
Misc: LMC 5956 86002740 3101-2493F  
Quant Time: 861020 12:30  
Injected at: 861016 02:40

BTL421

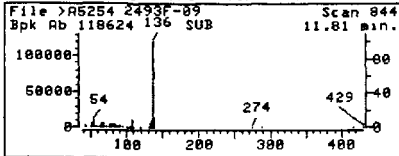
Compound No: 2  
Compound Name: 2-FLUOROPHENOL (SURROGATE)  
Scan Number: 241  
Retention Time: 5.59 min.  
Area: 135952  
Concentration: 86.13 UG/WIPE  
q-value: 100

220070

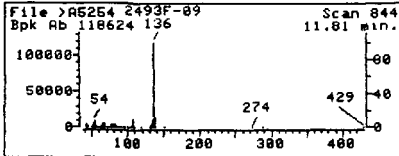
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

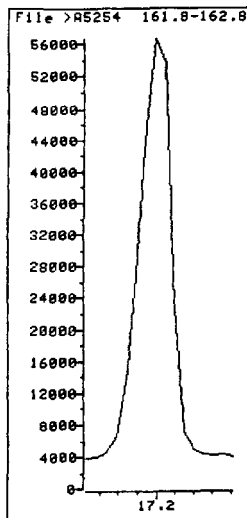
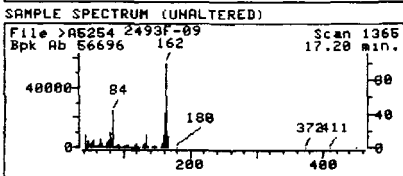
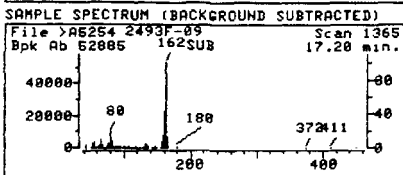
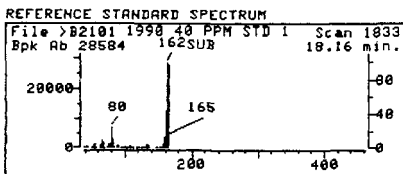


Data File: >A5254:E2  
Name: 2493F-09  
Misc: LMC 5996 86002740 3101-2493F  
Quant Time: 861020 12:30  
Injected at: 861016 02:40

BT1121

Compound No: 4 (ISTD)  
Compound Name: DG NAPHTHALENE  
Scan Number: 844  
Retention Time: 11.81 min.  
Area: 246434  
Concentration: 40.00 UG/WIPE  
q-value: 100

872010



Data File: >A5254::E2  
Name: 2493F-09  
Misc: LMC 5996 86002740 3101-2493F  
Quant Time: 861020 12:30  
Injected at: 861016 02:40

BTL021

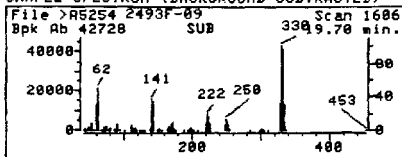
Compound No: 7 (ISTD)  
Compound Name: D10 ACENAPHTHENE  
Scan Number: 1365  
Retention Time: 17.20 min.  
Area: 129711  
Concentration: 40.00 UG/WIPE  
q-value: 100

672070

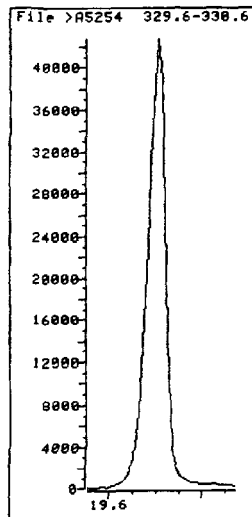
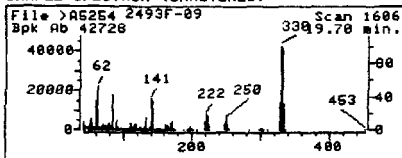
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



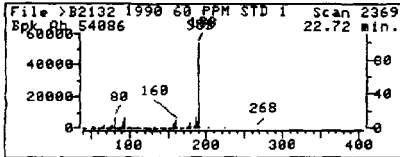
Data File: >A5254::E2  
Name: 2493F-09  
Misc: LMC 5996 86002740 3101-2493F  
Quant Time: 861020 12:30  
Injected at: 861016 02:40

BTL#21

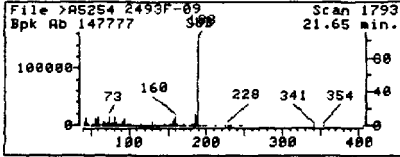
Compound No: 8  
Compound Name: 2,4,6-TRIBROMOPHENOL (SUFR)  
Scan Number: 1606  
Retention Time: 19.70 min.  
Area: 119123  
Concentration: 198.19 UG/WIFE  
q-value: 100

010000

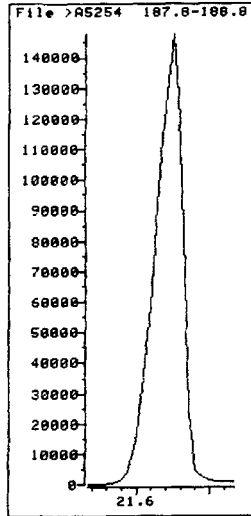
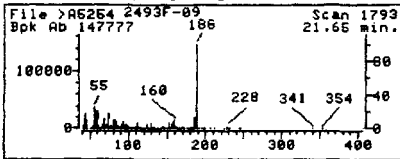
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >A5254:E2  
Name: 2493F-09  
Misc: LMC 5996 86002740 3101-2493F  
Quant Time: 861020 12:30  
Injected at: 861016 02:40

BT1021

Compound No: 11 (ISTD)  
Compound Name: D10 PHENANTHRENE  
Scan Number: 1793  
Retention Time: 21.65 min.  
Area: 389674  
Concentration: 40.00 UG/WIPE  
q-value: 100

102010



## Organics Analysis Data Sheet

(Page 1)

Laboratory Name: ENVIRONMENTAL ENGINEERS Case No: 3101-2493F  
 Lab Sample ID No: 86002748 QC Report No: \_\_\_\_\_  
 Sample Matrix: WIPES Contract No: SAS 3101-2493F  
 Data Release Authorized By: MARGARET C. WILKINSON Date Sample Received: 10/6/86

## Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: \_\_\_\_\_

Date Analyzed: \_\_\_\_\_

Conc/Dil Factor: \_\_\_\_\_ pH \_\_\_\_\_

Percent Moisture: (Not Decanted) \_\_\_\_\_

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	
67-64-1	Acetone	
75-15-0	Carbon Disulfide	
75-35-4	1, 1-Dichloroethene	
75-34-3	1, 1-Dichloroethane	
156-60-5	Trans-1, 2-Dichloroethene	
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	

CAS Number		ug/l or ug/Kg (Circle One)
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	
108-10-1	4-Methyl-2-Pentanone	
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	
108-90-7	Chlorobenzene	
100-41-4	Ethylbenzene	
100-42-5	Styrene	
	Total Xylenes	

0103010

## Data Reporting Qualifiers

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

- V** If the result is a value greater than or equal to the detection limit, report the value.
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample.
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10J). If limit of detection is 10 µg/l and a concentration of 3 µg/l is calculated, report as 3J.
- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ng/l in the final extract should be confirmed by GC/MS.
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Laboratory Name: ENVIRONMENTAL  
 Case No: 249

Sample Number  
2493F-10

Organics Analysis Data Sheet  
 (Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 10/6/86  
 Date Analyzed: 10/16/86  
 Conc/Dil Factor: 1  
 Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

*ug/wipe*  
*ug/ft<sup>2</sup> or ug/Kg*  
(Circle One)

CAS Number	Compound	Concentration
108-95-2	Phenol	
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	100
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	
120-83-2	2, 4-Dichlorophenol	100
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	100
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2, 4, 6-Trichlorophenol	100
95-95-4	2, 4, 5-Trichlorophenol	0.75
91-58-7	2-Chloronaphthalene	
88-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	

*ug/wipe*  
*ug/ft<sup>2</sup> or ug/Kg*  
(Circle One)

CAS Number	Compound	Concentration
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	
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	100
85-01-8	Phenanthrene	
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
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	
218-01-9	Chrysene	
117-84-0	Di-n-Octyl Phthalate	
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	Dibenz(a, h)Anthracene	
191-24-2	Benzo(g, h, i)Perylene	

(1)-Cannot be separated from diphenylamine

010388

*Handwritten notes and signatures at the bottom of the page.*

Laboratory Name: ENVIRODYNE  
Case No: 24931

Sample Number  
2493F-10

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.	no chlorophenols found			
2.				
3.				
4.				
5.				
6.				
7.				
8.				
9.				
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30.				

010334

QUANT REPORT

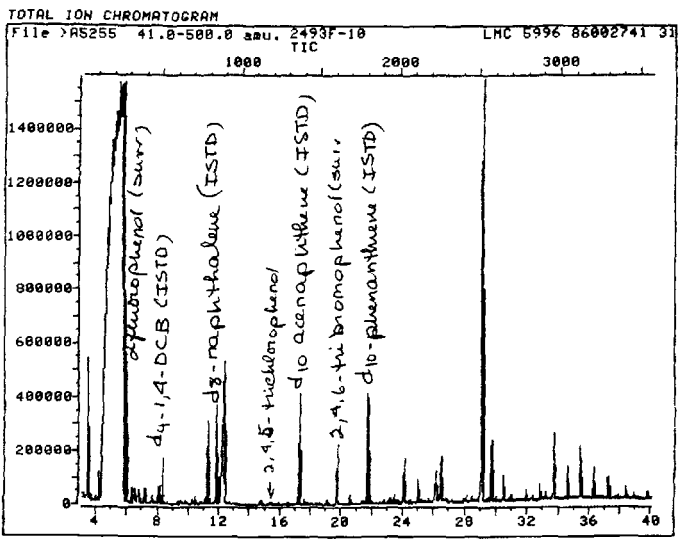
Operator ID: USER8      Quant Rev: 4      Quant Time: 861020 12:31  
 Output File: \*A5255::L2      Injected at: 861016 03:28  
 Data File: >A5255::E2      Dilution Factor: 1.000  
 Name: 2493F-10  
 Misc: LMC 5996 86002741 3101-2493F      BTL#22

ID File: #A5248::D3  
 Title: ID FILE FOR SAS 3101-2493F  
 Last Calibration: 861020 11:54

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *04 1,4-DICHLOROBENZENE	8.22	499	198255	40.00	UG/WIPE	93
2) 2-FLUOROPHENOL (SURROGATE)	5.92	276	211494	73.74	UG/WIPE	100
4) *08 NAPHTHALENE	11.81	848	468655	40.00	UG/WIPE	100
7) *010 ACENAPHTHENE	17.21	1370	231353	40.00	UG/WIPE	100
8) 2,4,6-TRIBROMOPHENOL (SURR)	19.68	1610	93979	87.66	UG/WIPE	100
10) 2,4,5-Trichlorophenol	15.28	1183	2082M	.68	UG/WIPE	63
11) *010 PHENANTHRENE	21.64	1800	425434	40.00	UG/WIPE	100

\* Compound is ISTD

010335  
 999010



010200

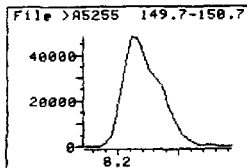
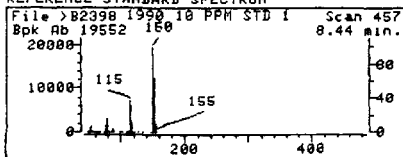
Data File: >AS255::E2  
Name: 2493F-10  
Misc: LMC 5996 86002741 3101-2493F

BTL022

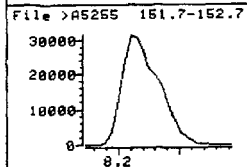
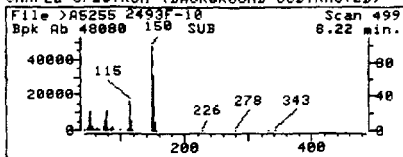
Id File: #AS240::D3  
Title: ID FILE FOR SRS 3101-2493F  
Last Calibration: 861020 11:54

Operator ID: USER0  
Quant Time: 861020 12:31  
Injected at: 861016 03:28

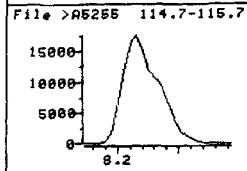
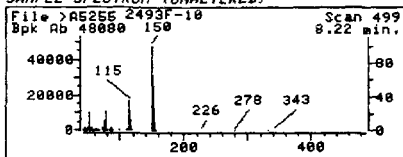
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



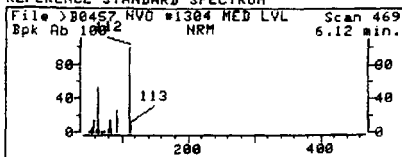
Data File: >A5255::E2  
Name: 2493F-10  
Misc: LMC 5996 B6002741 3101-2493F  
Quant Time: 861020 12:31  
Injected at: 861016 03:28

BTL#22

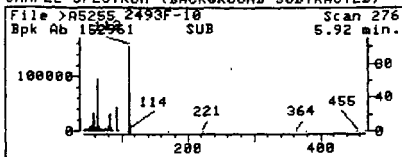
Compound No: 1 (ISTD)  
Compound Name: D4 1,4-DICHLOROBENZENE  
Scan Number: 499  
Retention Time: 8.22 min.  
Area: 198255  
Concentration: 40.00 UG/WIPE  
q-value: 93

010007

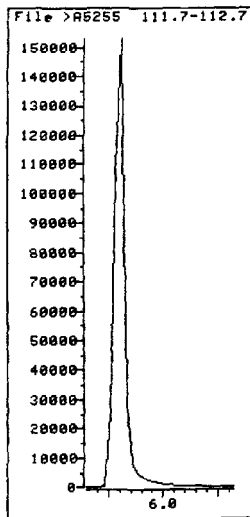
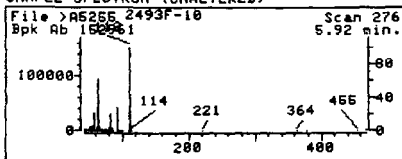
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



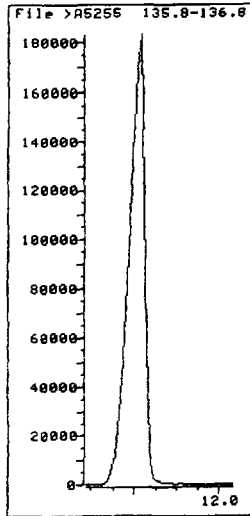
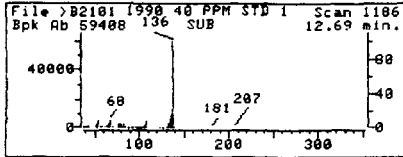
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Misc: LMC 5996 86002741 3101-2493F  
Quant Time: 861020 12:31  
Injected at: 861016 03:28

BTL#22

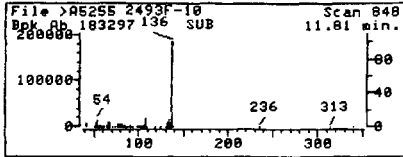
Compound No: 2  
Compound Name: 2-FLUOROPHENOL (SURROGATE)  
Scan Number: 276  
Retention Time: 5.92 min.  
Area: 211494  
Concentration: 73.74 UG/WIPE  
q-value: 100

010008

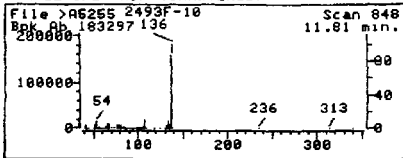
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >A5255::E2  
Name: 2493F-10  
Misc: LMC 5996 86002741 3101-2493F  
Quant Time: 861020 12:31  
Injected at: 861016 03:28

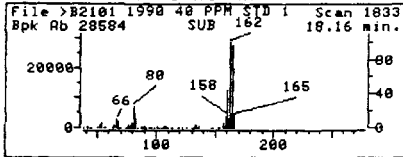
BTL#22

Compound No: 4 (ISTD)  
Compound Name: DG NAPHTHALENE  
Scan Number: 848  
Retention Time: 11.81 min.  
Area: 468655  
Concentration: 40.00 UG/WIPE  
q-value: 100

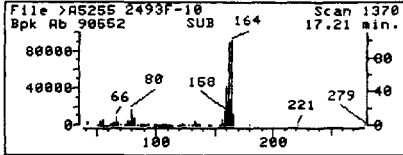
66010



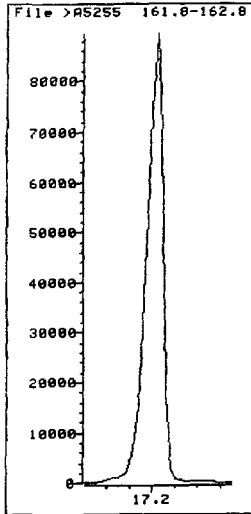
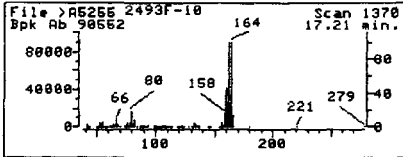
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

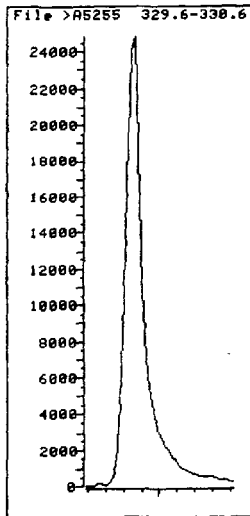
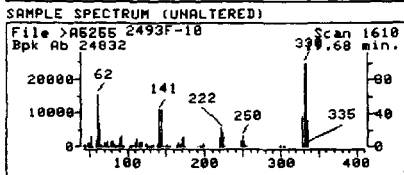
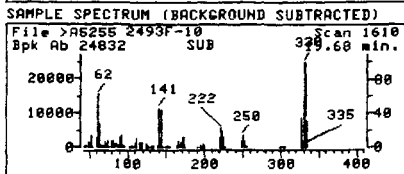
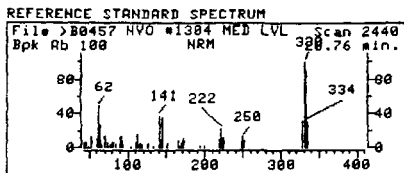


Data File: >A5255:E2  
Name: 2493F-10  
Misc: LMC 5996 86002741 3101-2493F  
Quant Time: 861020 12:31  
Injected at: 861016 03:28

BTL122

Compound No: 7 (ISTD)  
Compound Name: D10 ACENAPHTHENE  
Scan Number: 1370  
Retention Time: 17.21 min.  
Area: 231353  
Concentration: 40.00 UG/WIPE  
q-value: 100

010000

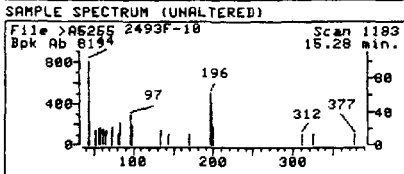
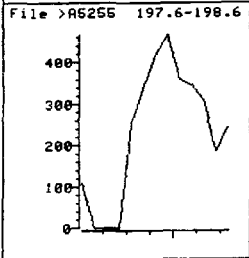
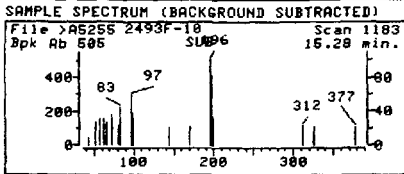
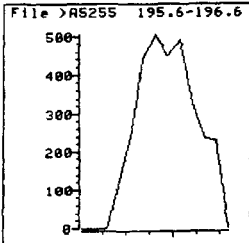
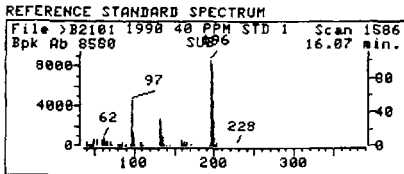


Data File: >A6255::E2  
Name: 2493F-10  
Misc: LMC 5996 86002741 3101-2493F  
Quant Time: 861020 12:31  
Injected at: 861016 03:28

BTL#22

Compound No: 8  
Compound Name: 2,4,6-TRIBROMOPHENOL (SURRE)  
Scan Number: 1610  
Retention Time: 19.68 min.  
Area: 93979  
Concentration: 87.66 UG/WIPE  
q-value: 100

100010



Data File: >A5255:E2  
Name: 2493F-10  
Misc: LMC 5996 86002741 3101-2493F  
Quant Time: 861020 12:31  
Injected at: 861016 03:28

BTL122

Compound No: 10  
Compound Name: 2,4,5-Trichlorophenol  
Scan Number: 1183  
Retention Time: 15.28 min.  
Area: 1891  
Concentration: .62 UG/WIPE  
q-value: 63

010092

Organics Analysis Data Sheet  
(Page 1)

Laboratory Name: ENVIRONMENTAL ENGRS Case No: 3101-2493F  
 Lab Sample ID No: 8600294 QC Report No: \_\_\_\_\_  
 Sample Matrix: WIDE Contract No: SAS 3101-2493F  
 Data Release Authorized By: Margaret C. Winters Date Sample Received: 10/6/86

## Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: \_\_\_\_\_

Date Analyzed: \_\_\_\_\_

Conc/Dil Factor: \_\_\_\_\_ pH \_\_\_\_\_

Percent Moisture: (Not Decanted) \_\_\_\_\_

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	
67-64-1	Acetone	
75-15-0	Carbon Disulfide	
75-35-4	1, 1-Dichloroethane	
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	

CAS Number		ug/l or ug/Kg (Circle One)
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	
108-10-1	4-Methyl-2-Pentanone	
591-78-6	2-Hexanone	
127-18-4	Tetrachloroethene	
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	
108-90-7	Chlorobenzene	
100-41-4	Ethylbenzene	
100-42-5	Styrene	
	Total Xylenes	

## Data Reporting Qualifiers

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

- Value** If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnote should read: U-Compound was analyzed for but not detected. The number is the minimum attainable detection limit for the sample
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10U). If limit of detection is 10 µg/l and a concentration of 3 µg/l is calculated, report as 3J.
- C** The flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides ≥ 10 ng/ul on the final extract should be confirmed by GC/MS.
- B** This flag is used when the analyte is found on the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Laboratory Name: ENVIRODYNE

Case No: 2493

Sample Number

2493 F-11

Organics Analysis Data Sheet  
(Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 10/6/86  
 Date Analyzed: 10/16/86  
 Conc/Dil Factor: 1  
 Percent Moisture (Decanted) \_\_\_\_\_

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/wipe ug/100ug/kg (Circle One)
108-95-2	Phenol	
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	100
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	
120-83-2	2, 4-Dichlorophenol	100
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	100
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2, 4, 6-Trichlorophenol	100
95-95-4	2, 4, 5-Trichlorophenol	500
91-58-7	2-Chloronaphthalene	
88-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	

CAS Number		ug/wipe ug/100ug/kg (Circle One)
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	
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	
87-85-5	Pentachlorophenol	500
85-01-8	Phenanthrene	
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
129-00-0	Pyrene	
95-68-7	Butylbenzylphthalate	
91-94-1	3, 3'-Dichlorobenzidine	
56-55-3	Benzo(a)Anthracene	
117-81-7	bis(2-Ethylhexyl)Phthalate	
218-01-9	Chrysene	
117-84-0	Di-n-Octyl Phthalate	
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	Dibenz(a, h)Anthracene	
191-24-2	Benzo(g, h, i)Perylene	

(1)-Cannot be separated from diphenylamine

00010

Laboratory Name: ENVIRONMENTAL  
Case No: 2493F

Sample Number  
2493F-11

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.	no chlorophenols found			
2.				
3.				
4.				
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010895

QUANT REPORT

Operator ID: USER8                    Quant Rev: 4    Quant Time: 861020 12:33  
 Output File: ^A5256::L2                    Injected at: 861016 04:15  
 Data File: >A5256::E2                    Dilution Factor: 1.000  
 Name: 2493F-11  
 Misc: LMC 5996 86002742 3101-2493F                    BTL#23

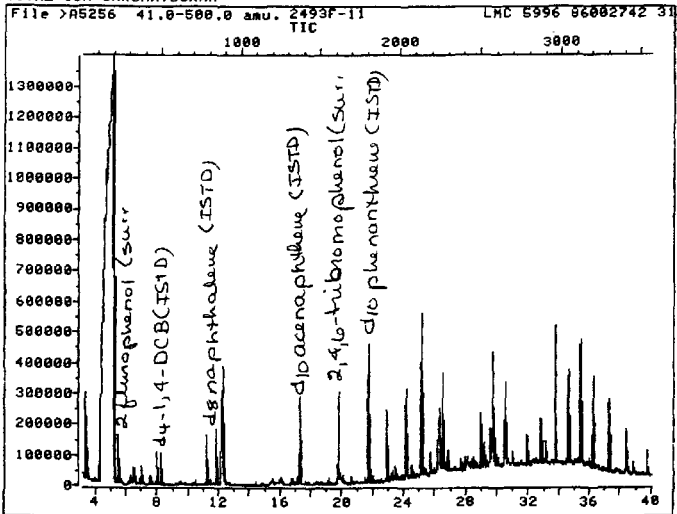
ID File: #A5248::D3  
 Title: ID FILE FOR SAS 3101-2493F  
 Last Calibration: 861020 11:54

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *04 1,4-DICHLOROBENZENE	8.16	491	69501	40.00	US/WIPE	91
2) 2-FLUOROPHENOL (SURROGATE)	9.46	229	76101	80.31	US/WIPE	100
4) *08 NAPHTHALENE	11.80	845	170306	40.00	US/WIPE	100
7) *010 ACENAPHTHENE	17.21	1370	144542	40.00	US/WIPE	100
8) 2,4,6-TRIBROMOPHENOL (SURR)	19.70	1612	139674	202.56	US/WIPE	100
11) *010 PHENANTHRENE	21.65	1801	516295	40.00	US/WIPE	100

\* Compound is ISTD

010200

TOTAL ION CHROMATOGRAM



Data File: >AS256::E2  
Name: 2493F-11  
Misc: LMC 5996 86002742 3101-2493F

BTL#23

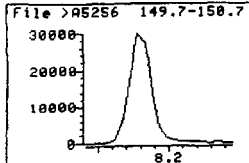
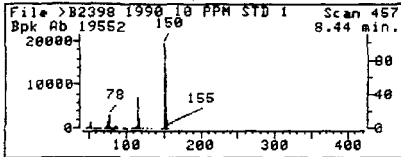
Id File: 0AS248::D3  
Title: ID FILE FOR SAS 3101-2493F  
Last Calibration: 861020 11:54

Operator ID: USEEB  
Quant Time: 861020 12:33  
Injected at: 861016 04:15

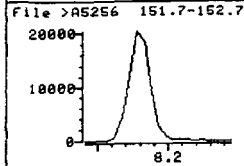
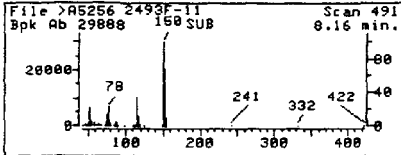
010007  
480010



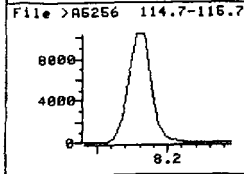
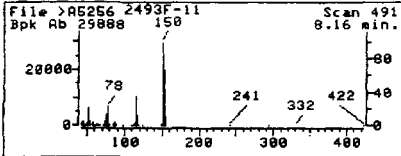
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



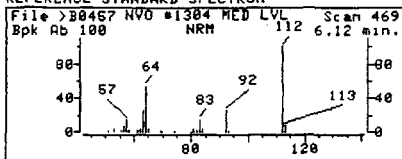
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Name: 2493F-11  
Misc: LMC 5996 86002742 3101-2493F  
Quant Time: 861020 12:33  
Injected at: 861016 04:15

BTL#23

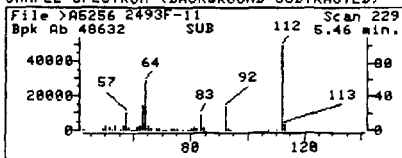
Compound No: 1 (ISTD)  
Compound Name: D4 1,4-DICHLOROBENZENE  
Scan Number: 491  
Retention Time: 8.16 min.  
Area: 65501  
Concentration: 40.00 UG/WIPE  
q-value: 91

010308

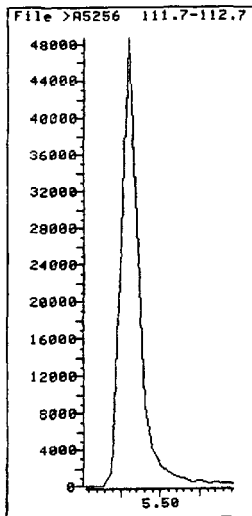
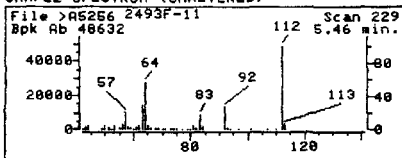
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



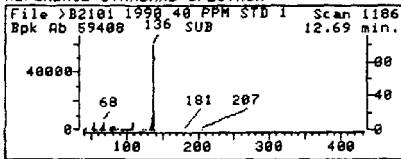
Data File: >A6256::E2  
Name: 2493F-11  
Misc: LMC 5996 86002742 3101-2493F  
Quant Time: 861020 12:33  
Injected at: 861016 04:15

BTL423

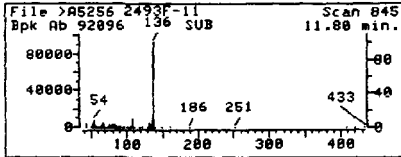
Compound No: 2  
Compound Name: 2-FLUOROPHENOL (SURROGATE)  
Scan Number: 229  
Retention Time: 5.46 min.  
Area: 76101  
Concentration: 80.31 US/WIPE  
q-value: 100

66010

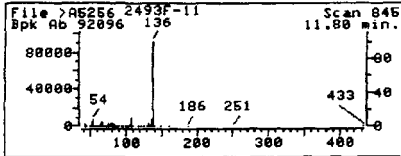
REFERENCE STANDARD SPECTRUM



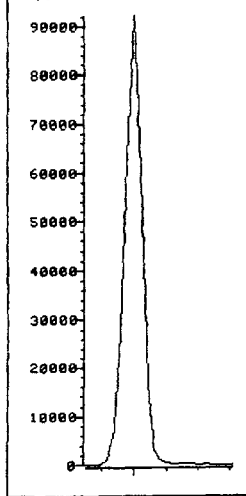
SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



File >A5256 135.8-136.8



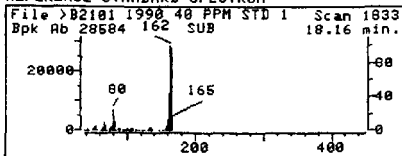
Data File: >A5256::E2  
Name: 2493F-11  
Misc: LMC 5996 86082742 3101-2493F  
Quant Time: 861020 12:33  
Injected at: 861016 04:15

BTL423

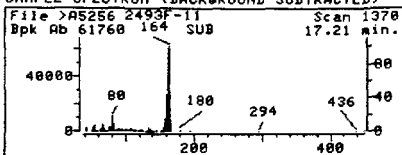
Compound No: 4 (ISTD)  
Compound Name: D8 NAPHTHALENE  
Scan Number: 845  
Retention Time: 11.80 min.  
Area: 170306  
Concentration: 40.00 UG/WIPE  
q-value: 100

010100

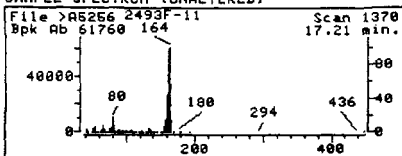
REFERENCE STANDARD SPECTRUM



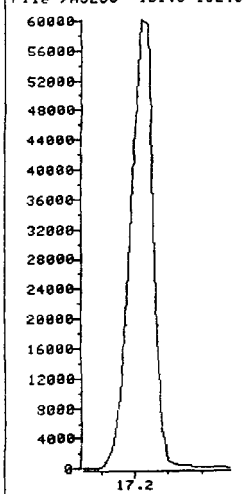
SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



File >A5256 161.8-162.8



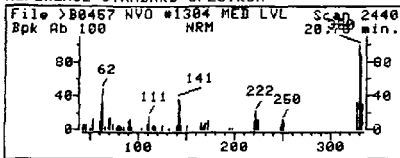
Data File: >A5256::E2  
Name: 2493F-11  
Misc: LMC 5996 86002742 3101-2493F  
Quant Time: 861020 12:33  
Injected at: 861016 04:15

BTL423

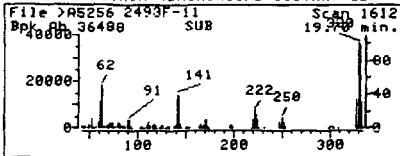
Compound No: 7 (ISTD)  
Compound Name: D10 ACENAPHTHENE  
Scan Number: 1370  
Retention Time: 17.21 min.  
Area: 144542  
Concentration: 40.00 UG/WIPE  
q-value: 100

010101

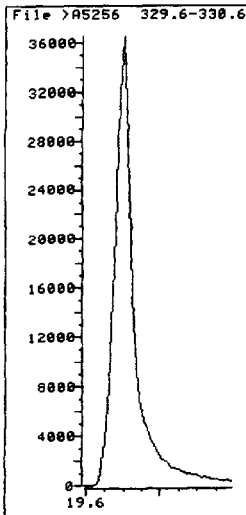
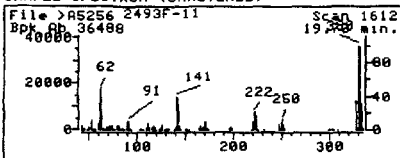
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)

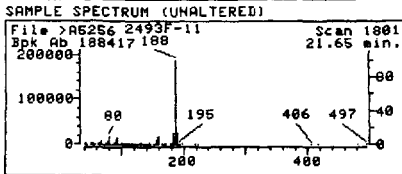
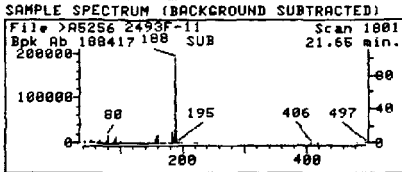
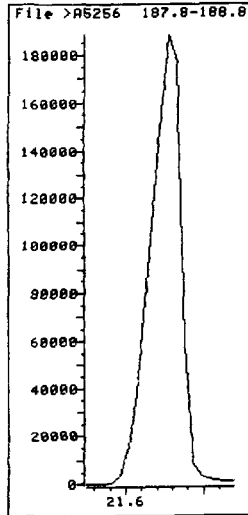
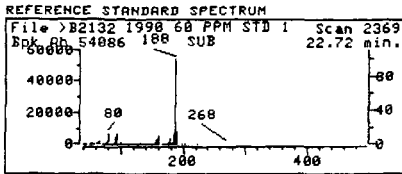


Data File: >AS256:E2  
Name: 2493F-11  
Misc: LMC 5996 86002742 3101-2493F  
Quant Time: 861020 12:33  
Injected at: 861016 04:15

BTL423

Compound No: 8  
Compound Name: 2,4,6-TRIBROMOPHENOL (SURR)  
Scan Number: 1612  
Retention Time: 19.70 min.  
Area: 135674  
Concentration: 202.56 UG/WIPE  
q-value: 100

001010



Data File: >A5256::E2  
 Name: 2493F-11  
 Misc: LMC 5996 86002742 3101-2493F  
 Quant Time: 061020 12:33  
 Injected at: 061016 04:15

BTL423

Compound No: 11 (ISTD)  
 Compound Name: D10 PHENANTHRENE  
 Scan Number: 1801  
 Retention Time: 21.65 min.  
 Area: 516295  
 Concentration: 40.00 UG/WIPE  
 q-value: 100

010103

Organics Analysis Data Sheet  
(Page 1)

Laboratory Name: ENVIRONMENTAL ENGRS 3101-2493F  
 Lab Sample ID No: 86002943 QC Report No: \_\_\_\_\_  
 Sample Matrix: WIDE Contact No: SAS 3101-2493F  
 Data Release Authorized By: Margaret C. White Date Sample Received: 10/6/86

## Volatile Compounds

Concentration: Low Medium (Circle One)

Date Extracted/Prepared: \_\_\_\_\_

Date Analyzed: \_\_\_\_\_

Conc/Dil Factor: \_\_\_\_\_ pH \_\_\_\_\_

Percent Moisture: (Not Decanted) \_\_\_\_\_

CAS Number		ug/l or ug/Kg (Circle One)
74-87-3	Chloromethane	
74-83-9	Bromomethane	
75-01-4	Vinyl Chloride	
75-00-3	Chloroethane	
75-09-2	Methylene Chloride	
67-64-1	Acetone	
75-15-0	Carbon Disulfide	
75-35-4	1, 1-Dichloroethane	
75-34-3	1, 1-Dichloroethane	
166-60-5	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	

CAS Number		ug/l or ug/Kg (Circle One)
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	
108-10-1	4-Methyl-2-Pentanone	
591-78-5	2-Hexanone	
127-18-4	Tetrachloroethene	
79-34-5	1, 1, 2, 2-Tetrachloroethane	
108-88-3	Toluene	
108-90-7	Chlorobenzene	
100-41-4	Ethylbenzene	
100-42-5	Styrene	
	Total Xylenes	

## Data Reporting Qualifiers

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

- Value** If the result is a value greater than or equal to the detection limit, report the value.
- U** Indicates compound was analyzed for but not detected. Report the minimum detection limit for the sample with the U (e.g., 10U) based on necessary concentration/dilution action. (This is not necessarily the instrument detection limit.) The footnotes should read: U-Compound was analyzed for but not detected. The number is the minimum ascertainable detection limit for the sample.
- J** Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed or when the mass spectral data indicated the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. (e.g., 10J). If limit of detection is 10 ug/l and a concentration of 3 ug/l is calculated, report as 2J.
- C** This flag applies to pesticide parameters where the identification has been confirmed by GC/MS. Single component pesticides  $\geq 10$  ng/l in the final extract should be confirmed by GC/MS.
- B** This flag is used when the analyte is found in the blank as well as a sample. It indicates possible/probable blank contamination and warns the data user to take appropriate action.
- Other** Other specific flags and footnotes may be required to properly define the results. If used, they must be fully described and such description attached to the data summary report.

Laboratory Name: ENVIRONMENTAL  
 Case No: 24931

Sample Number  
24931-12

Organics Analysis Data Sheet  
 (Page 2)

Semivolatile Compounds

Concentration: Low Medium (Circle One)  
 Date Extracted/Prepared: 10/6/86  
 Date Analyzed: 10-16-86  
 Conc/Dil Factor: 1  
 Percent Moisture (Decanted): \_\_\_\_\_

GPC Cleanup  Yes  No  
 Separatory Funnel Extraction  Yes  
 Continuous Liquid - Liquid Extraction  Yes

CAS Number		ug/wipe ug/100mg/Kg (Circle One)
108-95-2	Phenol	
111-44-4	bis(2-Chloroethyl)Ether	
95-57-8	2-Chlorophenol	10U
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	
62-72-1	Hexachloroethane	
98-95-3	Nitrobenzene	
78-59-1	Isophrone	
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	10U
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	10U
91-57-6	2-Methylnaphthalene	
77-47-4	Hexachlorocyclopentadiene	
88-06-2	2,4,6-Trichlorophenol	10U
95-95-4	2,4,5-Trichlorophenol	30U
91-58-7	2-Chloronaphthalene	
88-74-4	2-Nitroaniline	
131-11-3	Dimethyl Phthalate	
208-96-8	Acenaphthylene	
99-09-2	3-Nitroaniline	

CAS Number		ug/wipe ug/100mg/Kg (Circle One)
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	
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	50U
85-01-8	Phenanthrene	
120-12-7	Anthracene	
84-74-2	Di-n-Butylphthalate	
206-44-0	Fluoranthene	
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	
218-01-9	Chrysene	
117-84-0	Di-n-Octyl Phthalate	
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	Dibenz(a,h)Anthracene	
191-24-2	Benzo(g,h,i)Perylene	

(1)-Cannot be separated from diphenylamine

COFO



Laboratory Name: EN. RODYNE  
Case No: 2493F

Sample Number  
2493F-12

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.	no chlorophenols found			
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.				

010100

QUANT REPORT

Operator ID: USER8                      Quant Rev: 4    Quant Time: 861020 12:28  
 Output File: ^A5253::L2                      Injected at: 861016 01:53  
 Data File: >A5253::E2                      Dilution Factor: 1.000  
 Name: 2493F-12  
 Misc: LMC 5996 86002743 3101-2493F                      BTL#20

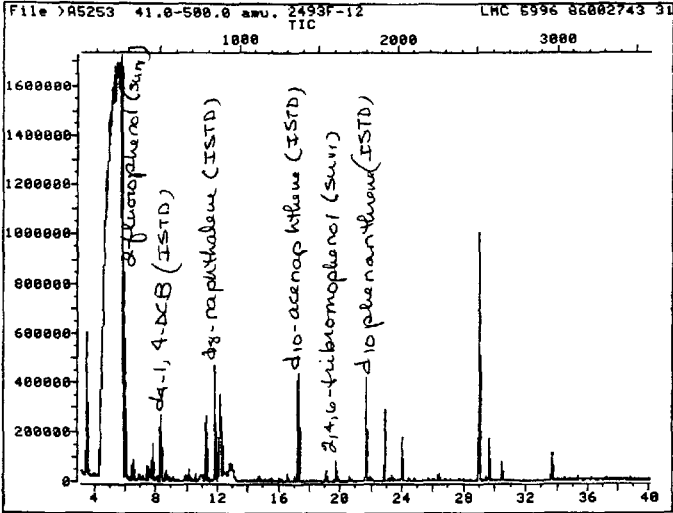
ID File: #A5248::03  
 Title: ID FILE FOR SAS 3101-2493F  
 Last Calibration: 861020 11:54

Compound	R.T.	Scan#	Area	Conc	Units	q
1) *D4 1,4-DICHLOROBENZENE	8.23	497	236463	40.00	UG/WIPE	94
2) 2-FLUOROPHENOL (SURROGATE)	5.97	278	215460	62.98	UG/WIPE	100
4) *D8 NAPHTHALENE	11.04	848	523115	40.00	UG/WIPE	100
7) *D10 ACENAPHTHENE	17.20	1368	235600	40.00	UG/WIPE	100
8) 2,4,6-TRIBROMOPHENOL (SURR)	19.65	1606	28148	25.78	UG/WIPE	100
11) *D10 PHENANTHRENE	21.63	1798	427473	40.00	UG/WIPE	100

\* Compound is ISTD

2010107

TOTAL ION CHROMATOGRAM



Data File: >A5253::E2  
Name: 2493F-12  
Misc: LMC 5996 86002743 3101-2493F

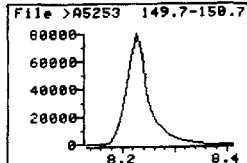
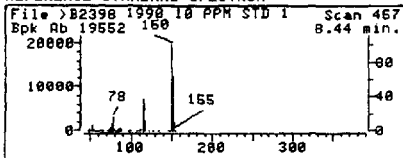
BTL420

Id File: #A5248::D3  
Title: ID FILE FOR SGS 3101-2493F  
Last Calibration: 861020 11:54

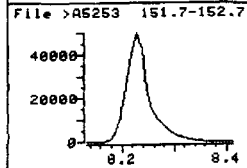
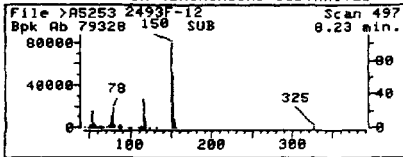
Operator ID: USER8  
Quant Time: 861020 12:28  
Injected at: 861016 01:53

010108

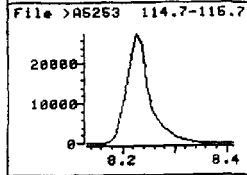
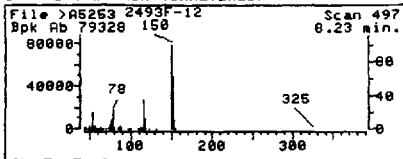
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



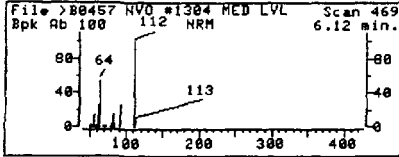
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 Misc: LHC 5996 86002743 3101-2493F  
 Quant Time: 861020 12:28  
 Injected at: 861016 01:53

BTL#20

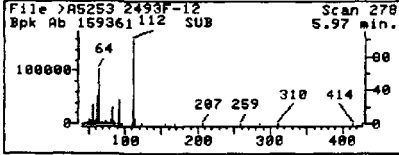
Compound No: 1 (ISTD)  
 Compound Name: D4 1,4-DICHLOROBENZENE  
 Scan Number: 497  
 Retention Time: 8.23 min.  
 Area: 236463  
 Concentration: 40.00 UG/WIPE  
 q-value: 94

601010

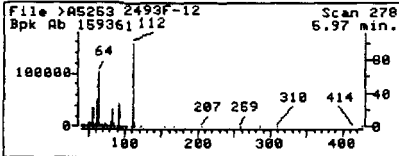
REFERENCE STANDARD SPECTRUM



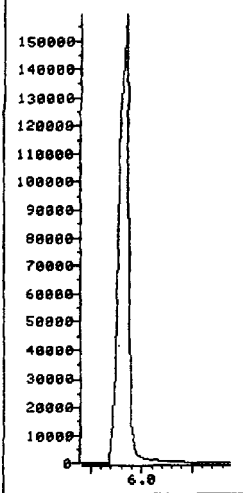
SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



File >A5253 111.7-112.7



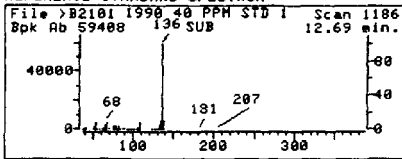
010110

Data File: >A5253::E2  
Name: 2493F-12  
Misc: LMC 5996 86002743 3101-2493F  
Quant Time: 861020 12:28  
Injected at: 861016 01:53

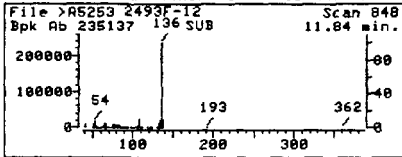
BTL020

Compound No: 2  
Compound Name: 2-FLUOROPHENOL (SURROGATE)  
Scan Number: 278  
Retention Time: 5.97 min.  
Area: 215460  
Concentration: 62.98 UG/WIPE  
q-value: 100

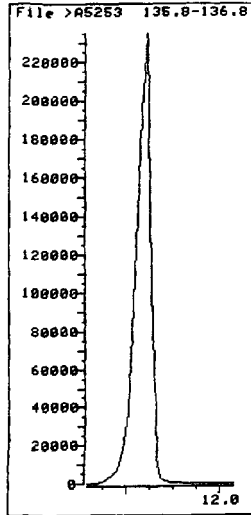
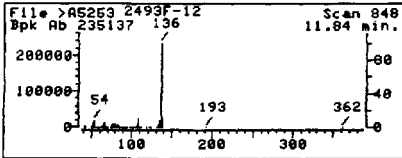
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



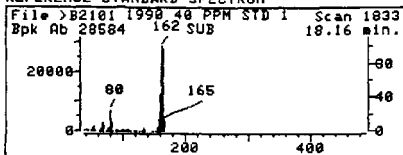
Data File: >A5253::E2  
Name: 2493F-12  
Misc: LMC 5996 86002743 3101-2493F  
Quant Time: 861020 12:28  
Injected at: 861016 01:53

BTL020

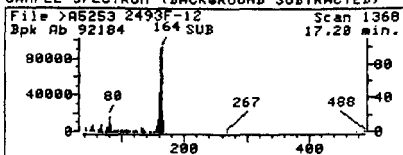
Compound No: 4 (ISTD)  
Compound Name: D8 NAPHTHALENE  
Scan Number: 848  
Retention Time: 11.84 min.  
Area: 523115  
Concentration: 40.00 UG/WIPE  
q-value: 100

010112

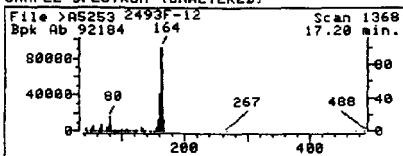
REFERENCE STANDARD SPECTRUM



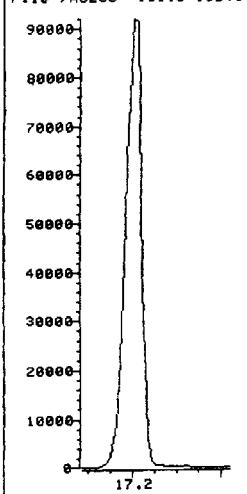
SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



File >A5253 161.8-162.8



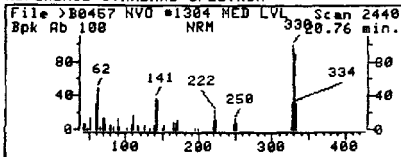
010112

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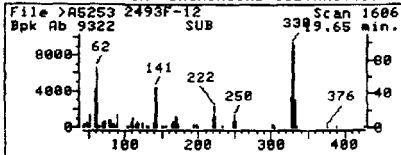
BTL#20

Compound No: 7 (ISTD)  
Compound Name: D10 ACENAPHTHENE  
Scan Number: 1368  
Retention Time: 17.20 min.  
Area: 235600  
Concentration: 40.00 UG/WIPE  
q-value: 100

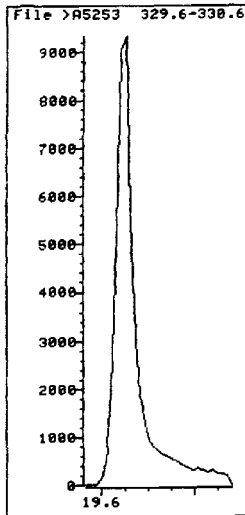
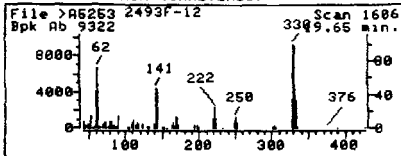
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >A5253::E2  
 Name: 2493F-12  
 Misc: LMC 5996 86002743 3101-2493F  
 Quant Time: 861020 12:28  
 Injected at: 861016 01:53

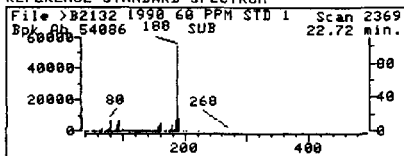
BTL120

Compound No: 0  
 Compound Name: 2,4,6-TRIBROMOPHENOL (SUER)  
 Scan Number: 1606  
 Retention Time: 19.65 min.  
 Area: 28140  
 Concentration: 25.78 UG/WIPE  
 q-value: 100

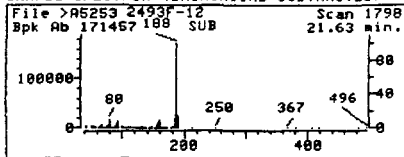
CT1010



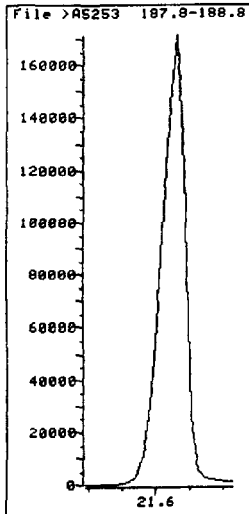
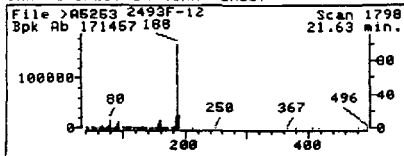
REFERENCE STANDARD SPECTRUM



SAMPLE SPECTRUM (BACKGROUND SUBTRACTED)



SAMPLE SPECTRUM (UNALTERED)



Data File: >A5253::E2  
Name: 2493F-12  
Misc: LMC 5996 86002743 3101-2493F  
Quant Time: 861020 12:28  
Injected at: 861016 01:53

BTL420

Compound No: 11 (ISTD)  
Compound Name: D10 PHENANTHRENE  
Scan Number: 1798  
Retention Time: 21.63 min.  
Area: 427473  
Concentration: 40.00 UG/UIPE  
q-value: 100

010414