User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 06/05/2002 Time: 08:43:41

Sample: AE12161
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 6/5/02 08:43 Ended: 6/5/02 08:43 Analyst: DLV
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		100		10.0

Comments for Sample AE12161 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 06-05-2002 Time: 08:44:08

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 4 of the 43 compounds in the LCS Standard were low.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052302\12161.D Vial: 14
Acq On : 23 May 2002 8:33 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 29 09:45:43 2002 Quant Results File: 052202.RES

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Wed May 29 09:44:24 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.90	152	3743484	40.00	ug/L	0.00
10) Napthalene-d8	13.39	136	14858131	40.00	ug/L	-0.01
17) Acenapthalene-d10	18.53	164	8209472	40.00	ug/L	0.00
29) Phenanthranene-d10	22.91	188	12389524	40.00	ug/L	0.00
37) Chrysene-d12	30.76	240	6789426	40.00	ug/L	-0.04
43) Perylene-d12	34.65	264	2858874	40.00	ug/L	-0.03
System Monitoring Compounds						
27) TCMX	20.50	209	1510001	39.86	ug/L	0.00
Spiked Amount 40.000			Recovery = 99.65%			
Target Compounds						Qvalue
2) n-nitros-dimethyl amine	0.00	74	0	N.D.	d	
3) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
4) 1,3-Dichlorobenzene	0.00	146	0	N.D.		
5) 1,4-Dichlorobenzene	0.00	146	0	N.D.		
6) 1,2-Dichlorobenzene	0.00	146	0	N.D.		
7) 2,2'oxybis(1-Chloropropane	0.00	45	0	N.D.		
8) N-nitroso-di-propylamine	0.00	70	0	N.D.		
9) Hexachloroethane	0.00	117	0	N.D.		
11) Nitrobenzene	0.00	77	0	N.D.		
12) Isophorone	0.00	82	0	N.D.		
13) bis(2-Chloroethoxy) methan	0.00	93	0	N.D.		
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
15) Napthalene	0.00	128	0	N.D.		
16) Hexachlorobutadiene	0.00	225	0	N.D.		
18) 2-Chloronaphthalene	0.00	162	0	N.D.		
19) Dimethylphthalate	0.00	163	0	N.D.		
20) 2,6-Dinitrotoluene	0.00	165	0	N.D.		
21) Acenapthylene	0.00	152	0	N.D.		
22) Acenapthalene	0.00	153	0	N.D.		
23) 2,4-Dinitrotoluene	0.00	165	0	N.D.		
24) Diethylphthalate	0.00	149	0	N.D.		
25) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.		
26) Fluorene	0.00	166	0	N.D.		
28) Azobenzene	20.50	77	26102	N.D.		
30) Nnitrosodiphenylamine	20.51	169	28619	N.D.		
31) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		
32) Hexachlorobenzene	0.00	284	0	N.D.		
33) Phenanthrene	0.00	178	0	N.D.		
34) Anthracene	0.00	178	0	N.D.		
35) Di-n-butylphthalate	0.00	149	0	N.D.		

(#) = qualifier out of range (m) = manual integration

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Title : 508 Modified GC/MS scan
Last Update : Wed May 29 09:44:24 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

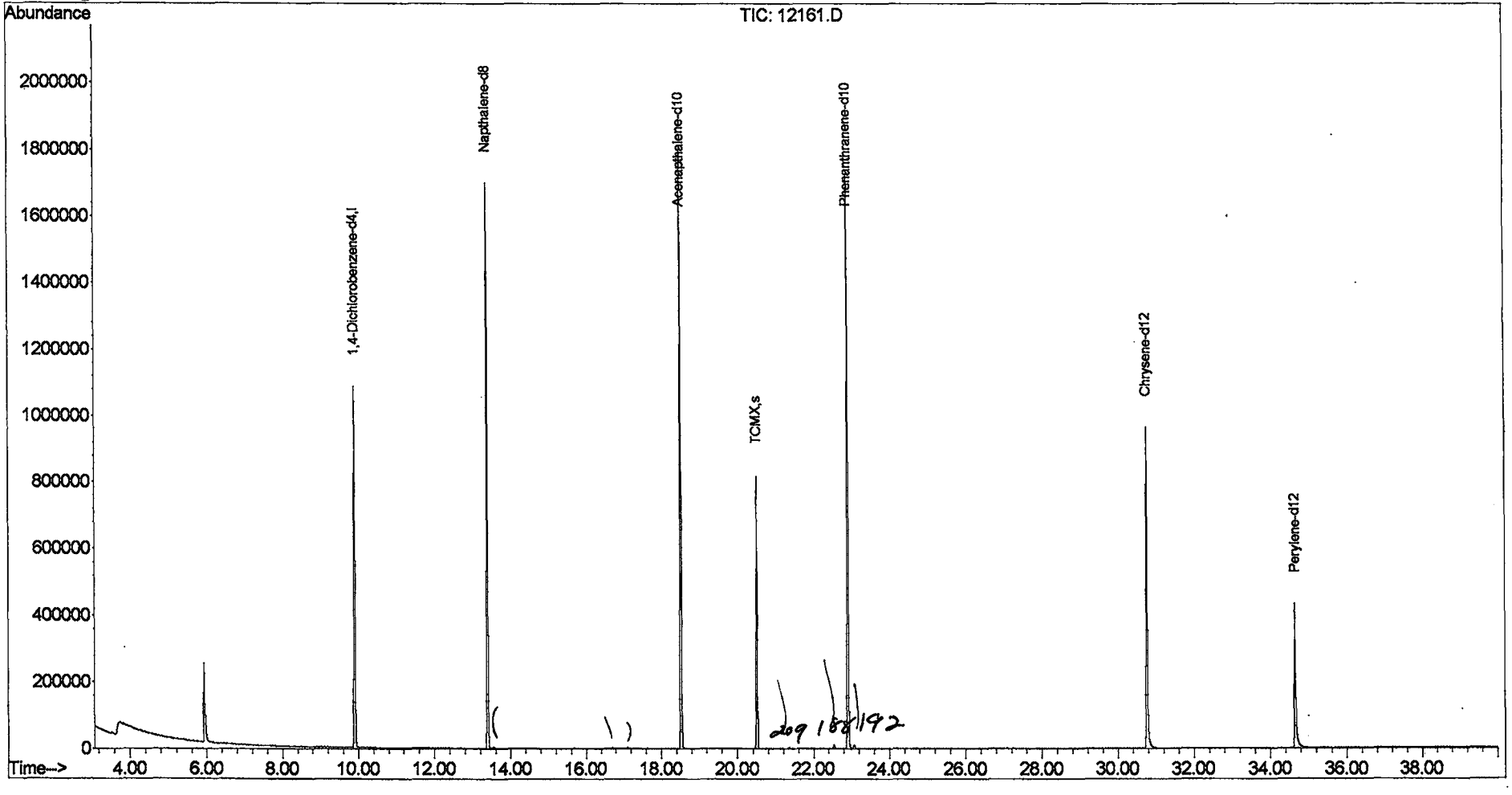
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
36) Fluoroanthene	0.00	202	0	N.D.	
38) Pyrene	0.00	202	0	N.D.	
39) Butylbenzylphthalate	0.00	149	0	N.D.	
40) Benzo(a)anthracene	30.76	228	14343	N.D.	
41) Bis(2-ethylhexyl)phthalate	0.00	149	0	N.D.	
42) Chrysene	0.00	228	0	N.D.	
44) Di-n-octylphthalate	0.00	149	0	N.D.	
45) Benzo(b)fluoranthene	0.00	252	0	N.D.	
46) Benzo(k)fluoranthene	0.00	252	0	N.D.	
47) Benzo(a)pyrene	0.00	252	0	N.D.	
48) Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
49) Dibenz(a,h)anthracene	0.00	278	0	N.D.	
50) Benzo(g,h,i)perylene	0.00	276	0	N.D.	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\052302\12161.D Vial: 14
Acq On : 23 May 2002 8:33 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 29 9:52 2002 Quant Results File: 052202.RES

Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Wed May 29 09:44:24 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\052302\12161.D
Acq On : 23 May 2002 8:33 pm
Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 14
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan

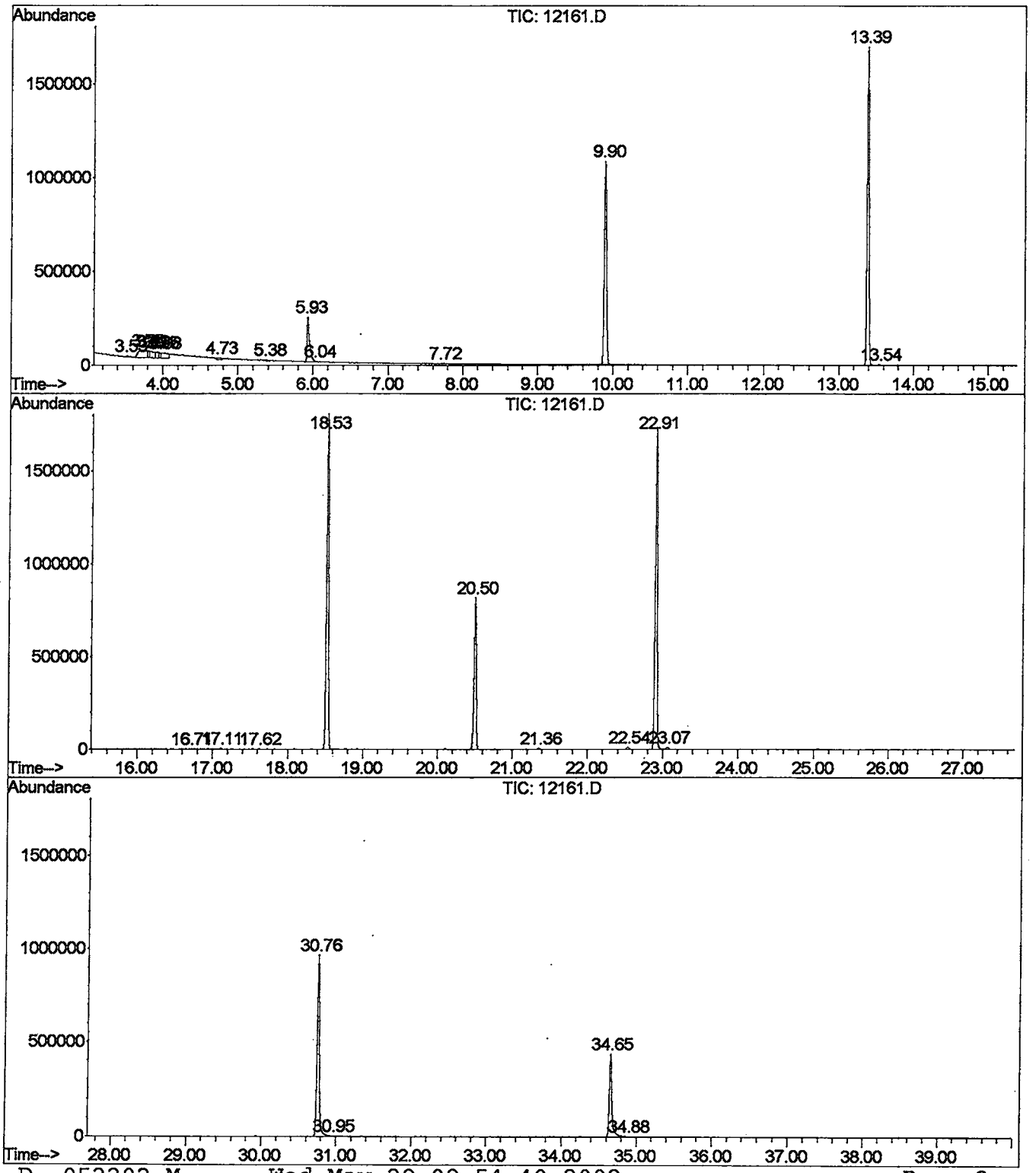
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.526	72	76	81	PV 7	4449	77614	0.23%	0.043%
2	3.743	94	113	122	PV 6	38416	2958381	8.77%	1.633%
3	3.808	122	124	127	VV 4	36471	550022	1.63%	0.304%
4	3.831	127	128	140	VV 6	33439	1448454	4.29%	0.800%
5	3.914	140	142	147	VV 5	30104	730513	2.16%	0.403%
6	3.949	147	148	152	VV 3	28317	444049	1.32%	0.245%
7	3.978	152	153	171	VV 9	27144	1696309	5.03%	0.937%
8	4.730	277	281	291	VV 7	8352	336827	1.00%	0.186%
9	5.377	385	391	401	VV 8	3974	160748	0.48%	0.089%
10	5.929	479	485	503	PV	233860	5455806	16.17%	3.012%
11	6.041	503	504	508	VV 3	3554	45517	0.13%	0.025%
12	7.715	786	789	795	PV 4	2635	58654	0.17%	0.032%
13	9.901	1152	1161	1183	PV	1076275	22453555	66.53%	12.397%
14	13.391	1742	1755	1773	PV	1668023	30319898	89.84%	16.740%
15	13.544	1777	1781	1792	VV 3	3865	71971	0.21%	0.040%
16	16.705	2315	2319	2326	PV 2	2966	48264	0.14%	0.027%
17	17.110	2383	2388	2395	VV 5	4123	62073	0.18%	0.034%
18	17.621	2471	2475	2479	VV 3	2083	24287	0.07%	0.013%
19	18.532	2605	2630	2647	PV	1780731	33648716	99.71%	18.578%
20	20.506	2954	2966	2974	BV	809853	14933624	44.25%	8.245%
21	21.358	3103	3111	3118	VV 5	4879	84967	0.25%	0.047%
22	22.539	3301	3312	3319	VV 2	11474	207181	0.61%	0.114%
23	22.909	3360	3375	3395	PV 2	1706061	33748117	100.00%	18.633%
24	23.068	3395	3402	3420	VV	10747	223480	0.66%	0.123%
25	30.759	4700	4711	4742	PV 2	953909	20760488	61.52%	11.462%
26	30.953	4742	4744	4757	VV 6	2450	54226	0.16%	0.030%
27	34.655	5364	5374	5410	PV	435309	10490712	31.09%	5.792%
28	34.878	5410	5412	5419	VV 5	1273	23492	0.07%	0.013%

Sum of corrected areas: 181117946

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\052302\12161.D
Operator : Mclean
Acquired : 23 May 2002 8:33 pm using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 5/22ad
Misc Info :
Vial Number: 14
Quant File :052202.RES (Chemstation Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\12161.D
Acq On : 23 May 2002 8:33 pm
Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 14
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

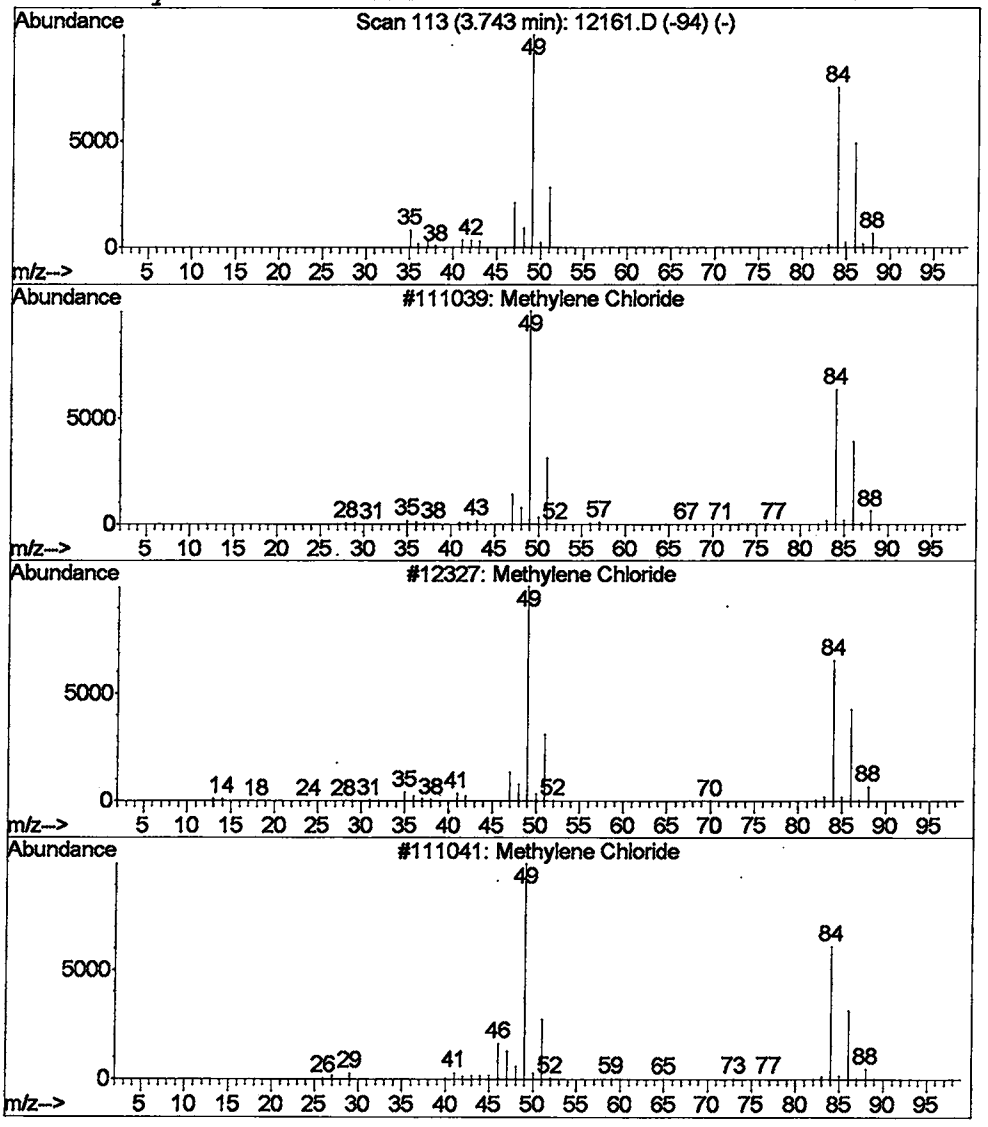
Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

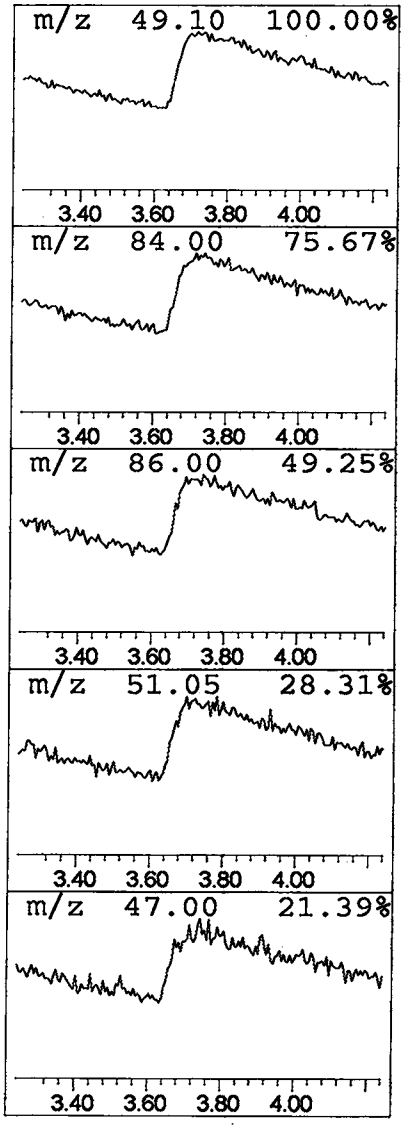
****

Peak Number 1 Methylene Chloride Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.75	5.27 ug/L	2958380	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methylene Chloride	84	CH2Cl2	000075-09-2	91
2	Methylene Chloride	84	CH2Cl2	000075-09-2	91
3	Methylene Chloride	84	CH2Cl2	000075-09-2	86
4	Methylene Chloride	84	CH2Cl2	000075-09-2	58





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\12161.D Vial: 14
Acq On : 23 May 2002 8:33 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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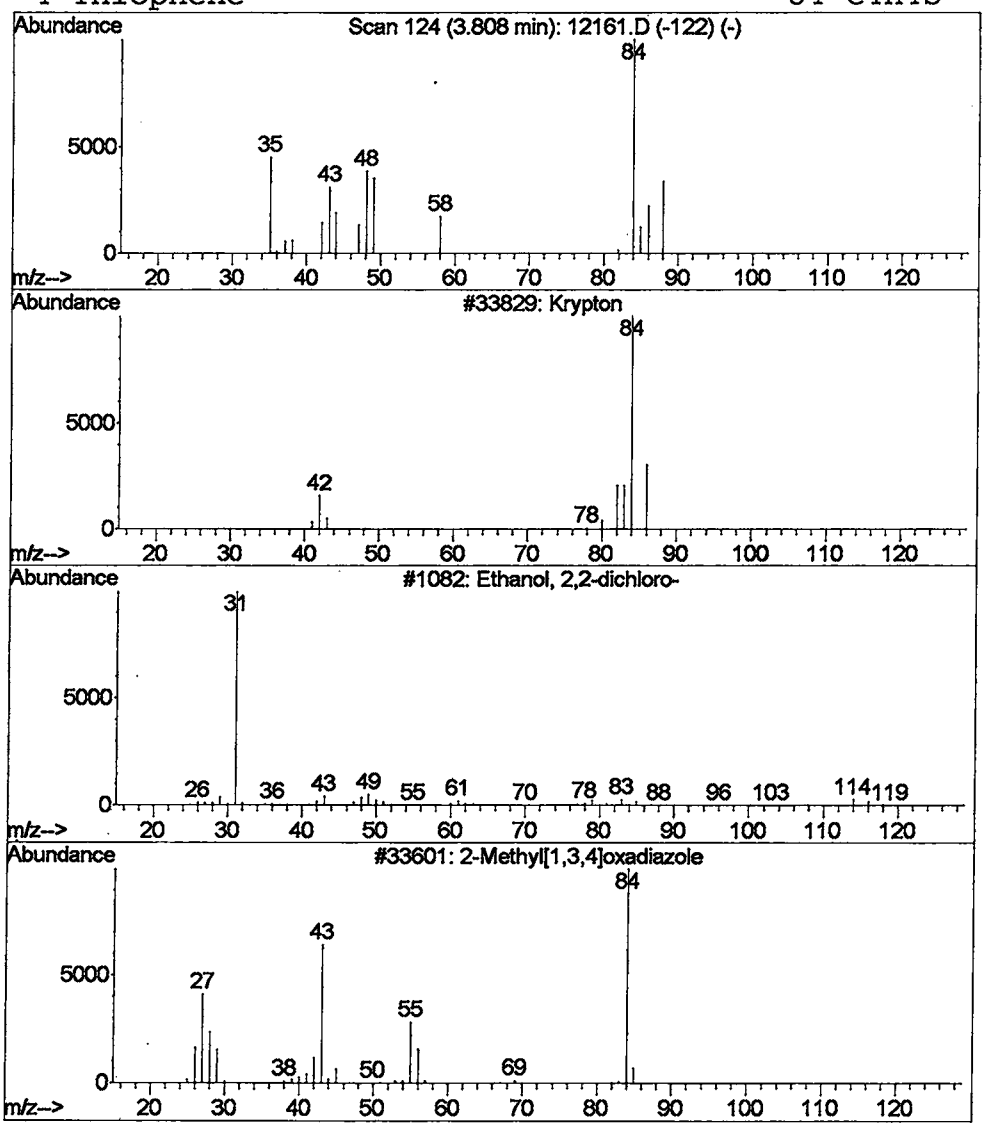
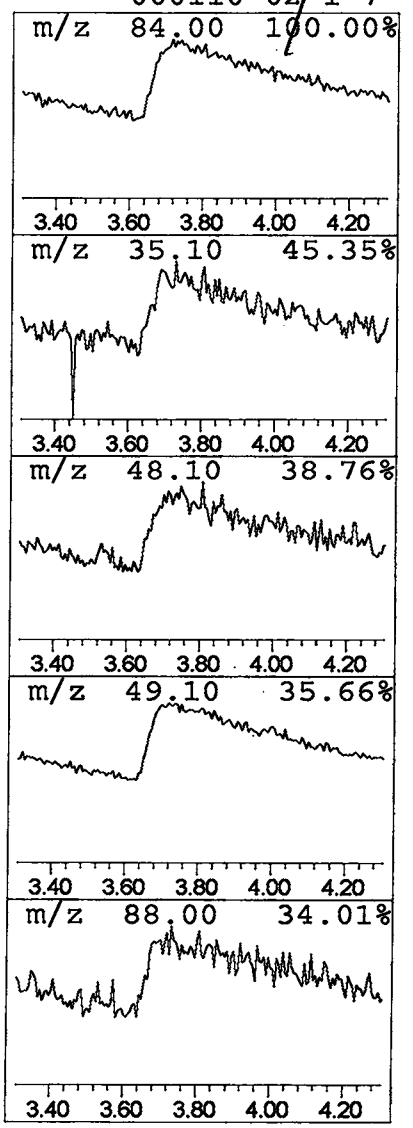
Peak Number 2 Krypton Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.81	0.98 ug/L	550022	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Krypton	84	Kr	007439-90-9	9
2	Ethanol, 2,2-dichloro-	114	C2H4Cl2O	000598-38-9	9
3	2-Methyl[1,3,4]oxadiazole	84	C3H4N2O	003451-51-2	7
4	Thiophene	84	C4H4S	000110-02-1	7





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\12161.D Vial: 14
Acq On : 23 May 2002 8:33 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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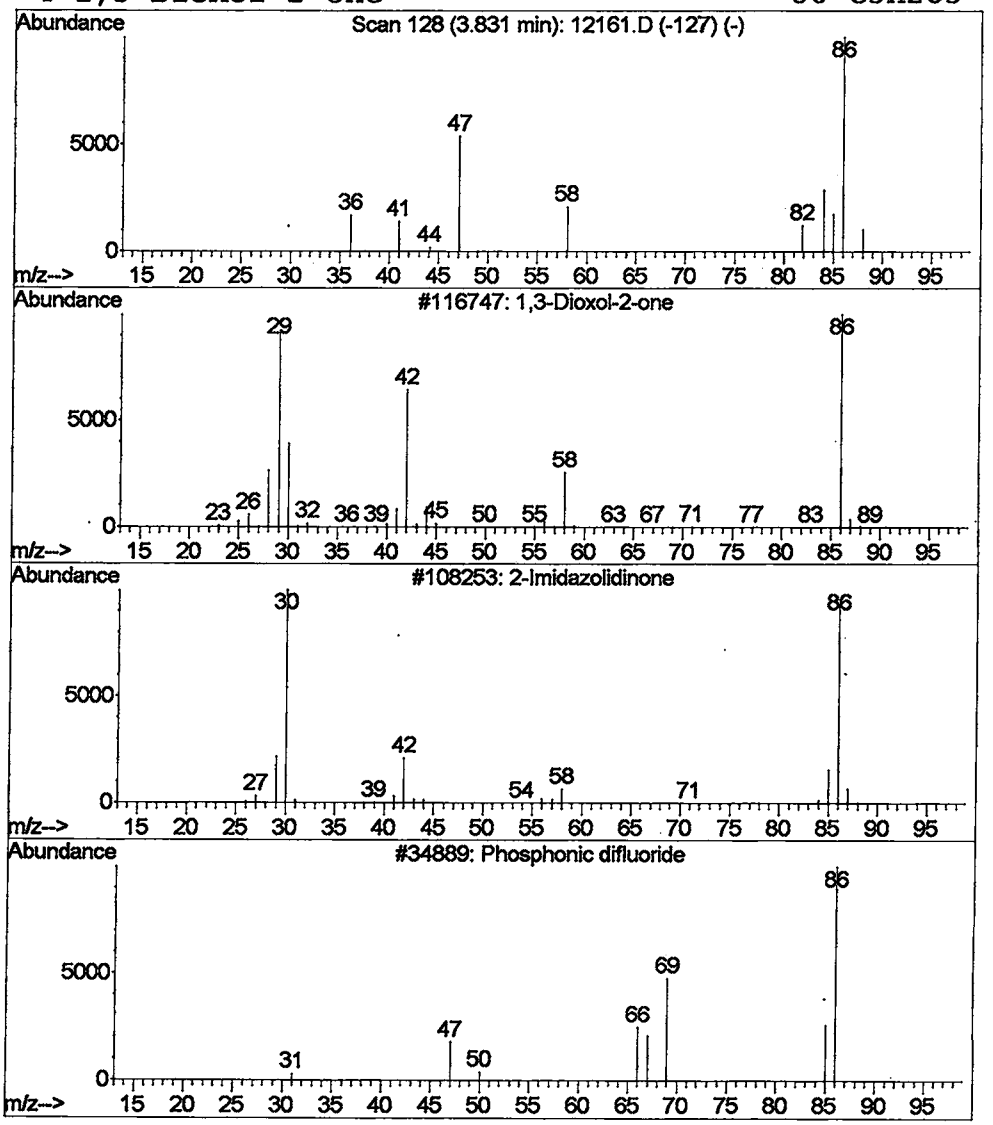
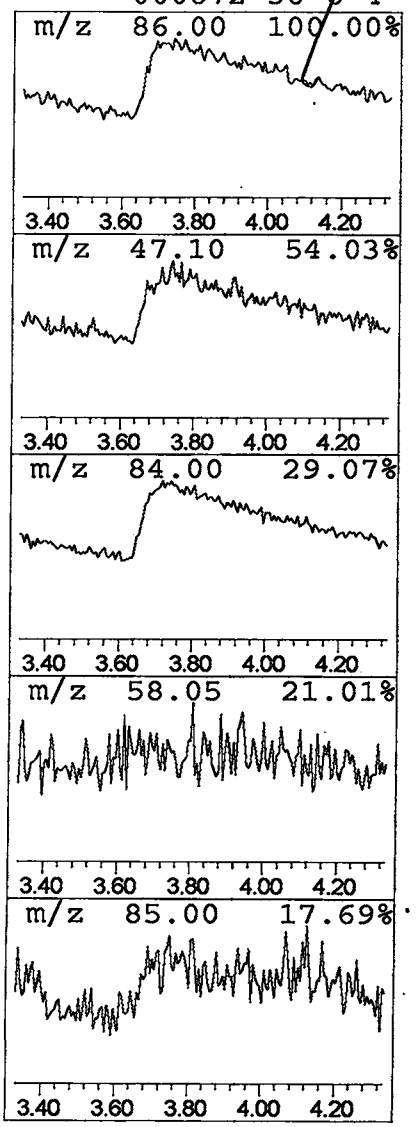
Peak Number 3 1,3-Dioxol-2-one Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.83	2.58 ug/L	1448450	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxol-2-one	86	C3H2O3	000872-36-6	7
2	2-Imidazolidinone	86	C3H6N2O	000120-93-4	7
3	Phosphonic difluoride	86	F2HOP	014939-34-5	5
4	1,3-Dioxol-2-one	86	C3H2O3	000872-36-6	4





Library Search Compound Report

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Acq On : 23 May 2002 8:33 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
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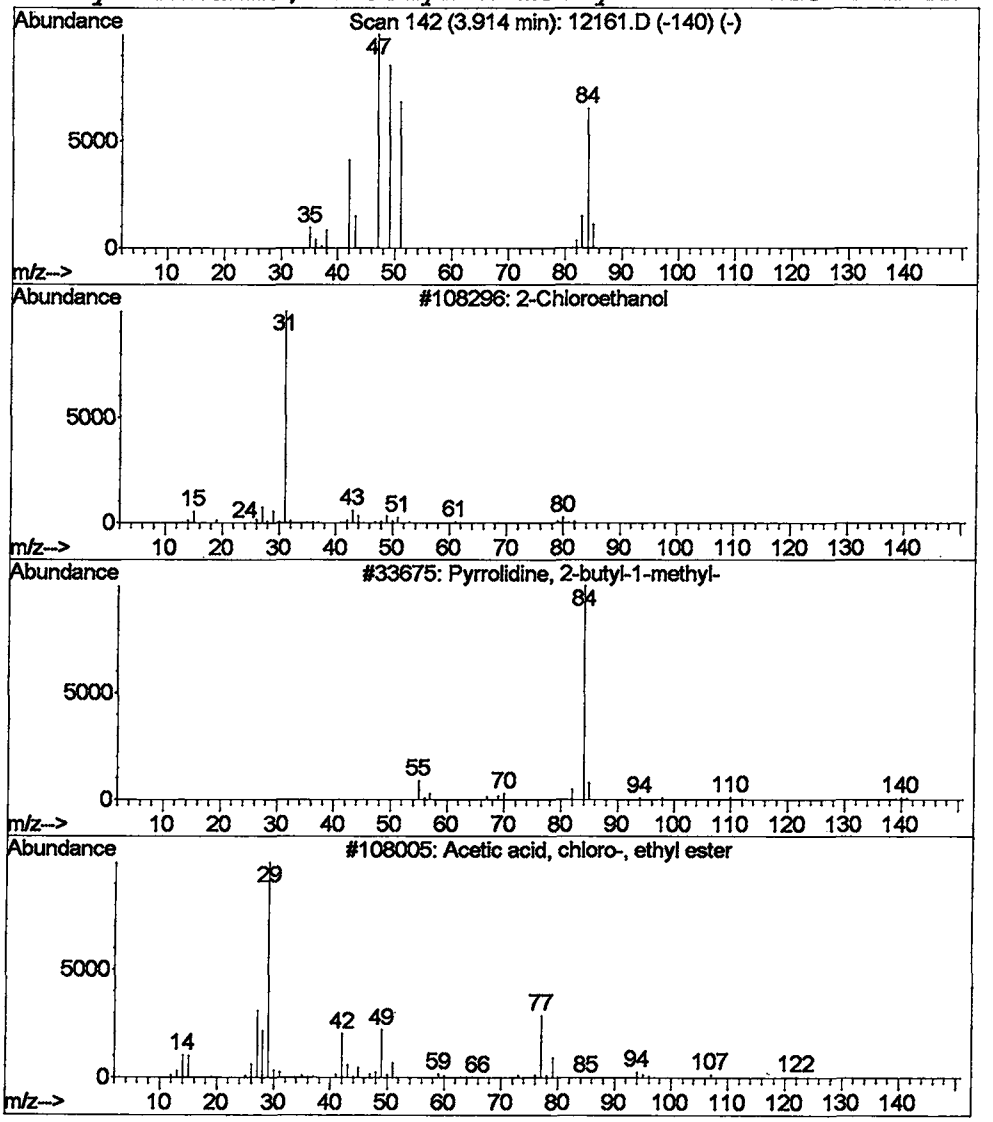
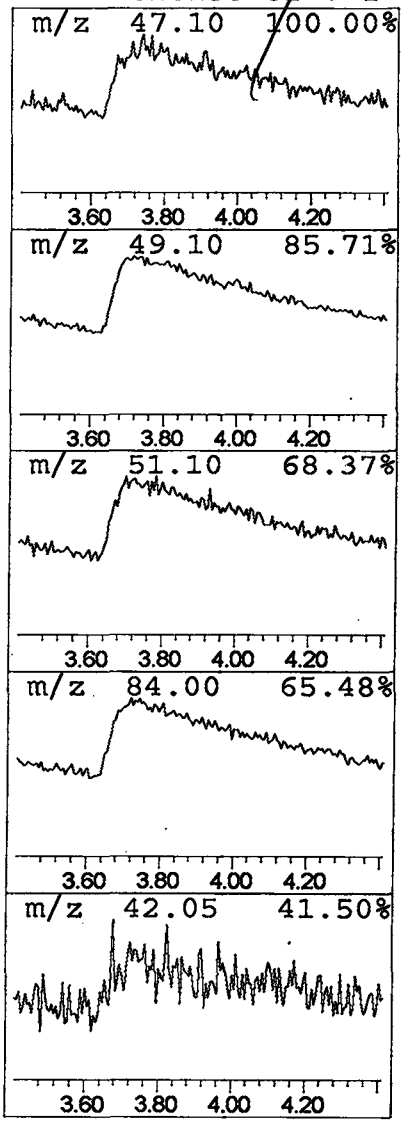
Peak Number 4 2-Chloroethanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.91	1.30 ug/L	730513	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloroethanol	80	C2H5ClO	000107-07-3	2
2	Pyrrolidine, 2-butyl-1-methyl-	141	C9H19N	003447-03-8	2
3	Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	2
4	Pyrrolidine, 2-ethyl-1-methyl-	113	C7H15N	026158-82-7	1





Library Search Compound Report

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Acq On : 23 May 2002 8:33 pm
Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 14
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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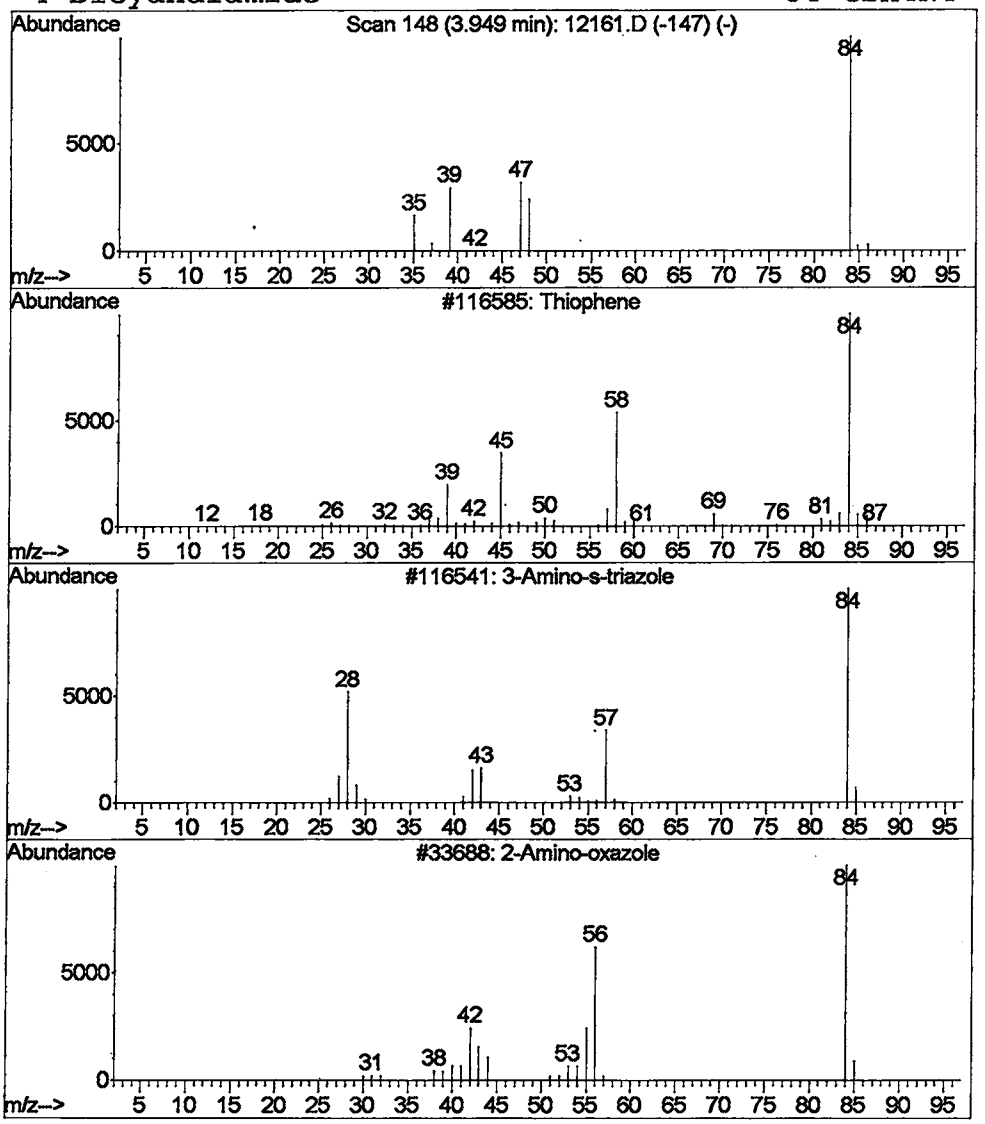
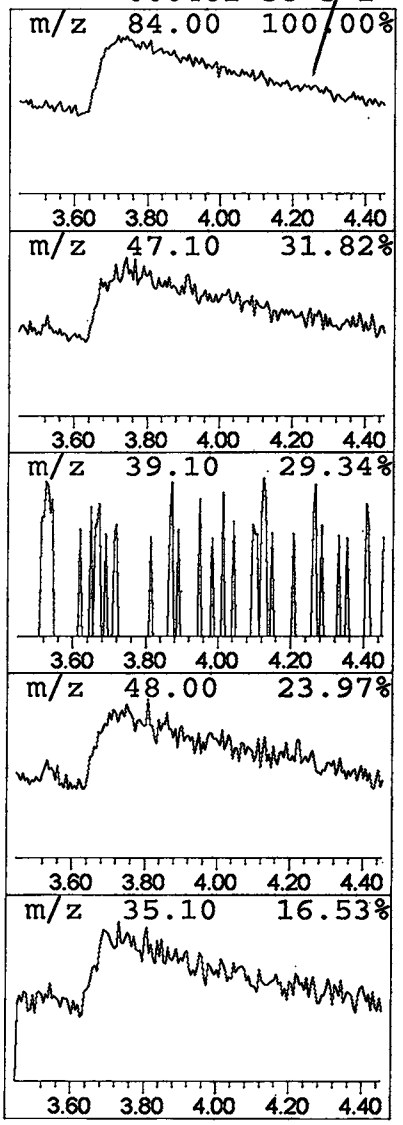
Peak Number 5 Thiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.95	0.79 ug/L	444049	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thiophene	84	C4H4S	000110-02-1	4
2	3-Amino-s-triazole	84	C2H4N4	000061-82-5	2
3	2-Amino-oxazole	84	C3H4N2O	1000144-64-0	2
4	Dicyandiamide	84	C2H4N4	000461-58-5	2





Library Search Compound Report

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Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 14
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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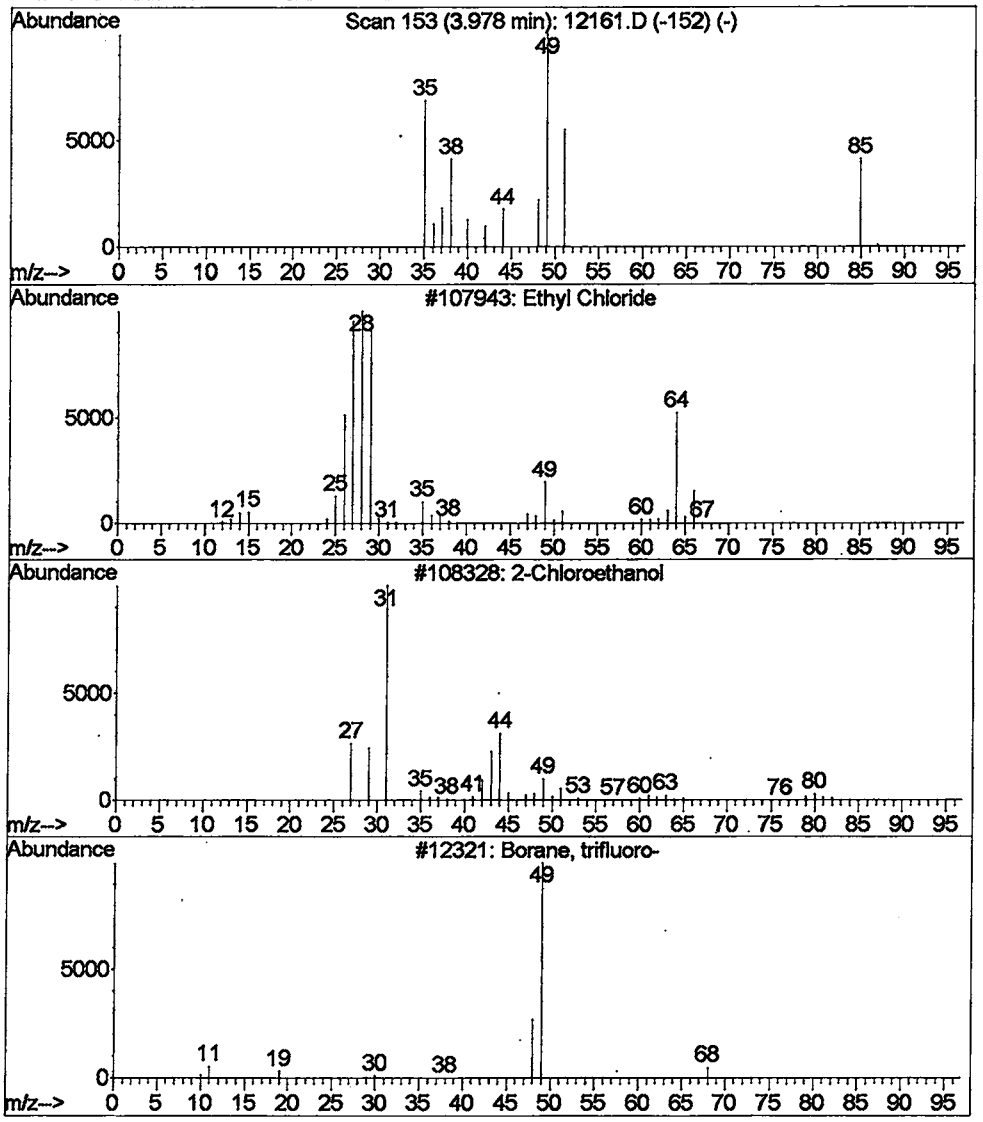
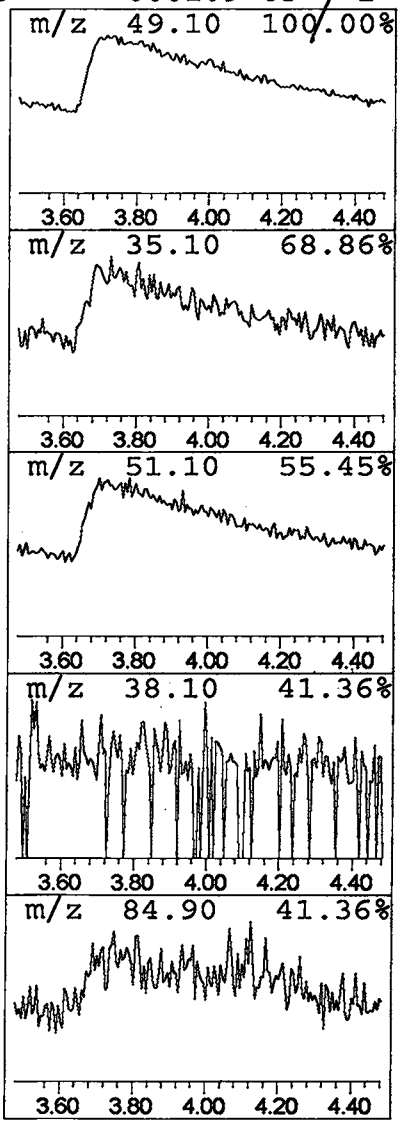
Peak Number 6 Ethyl Chloride Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.98	3.02 ug/L	1696310	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl Chloride	64	C2H5Cl	000075-00-3	36
2	2-Chloroethanol	80	C2H5ClO	000107-07-3	2
3	Borane, trifluoro-	68	BF3	007637-07-2	2
4	Boron trifluoride etherate	142	C4H10BF3O	000109-63-7	2





Library Search Compound Report

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Acq On : 23 May 2002 8:33 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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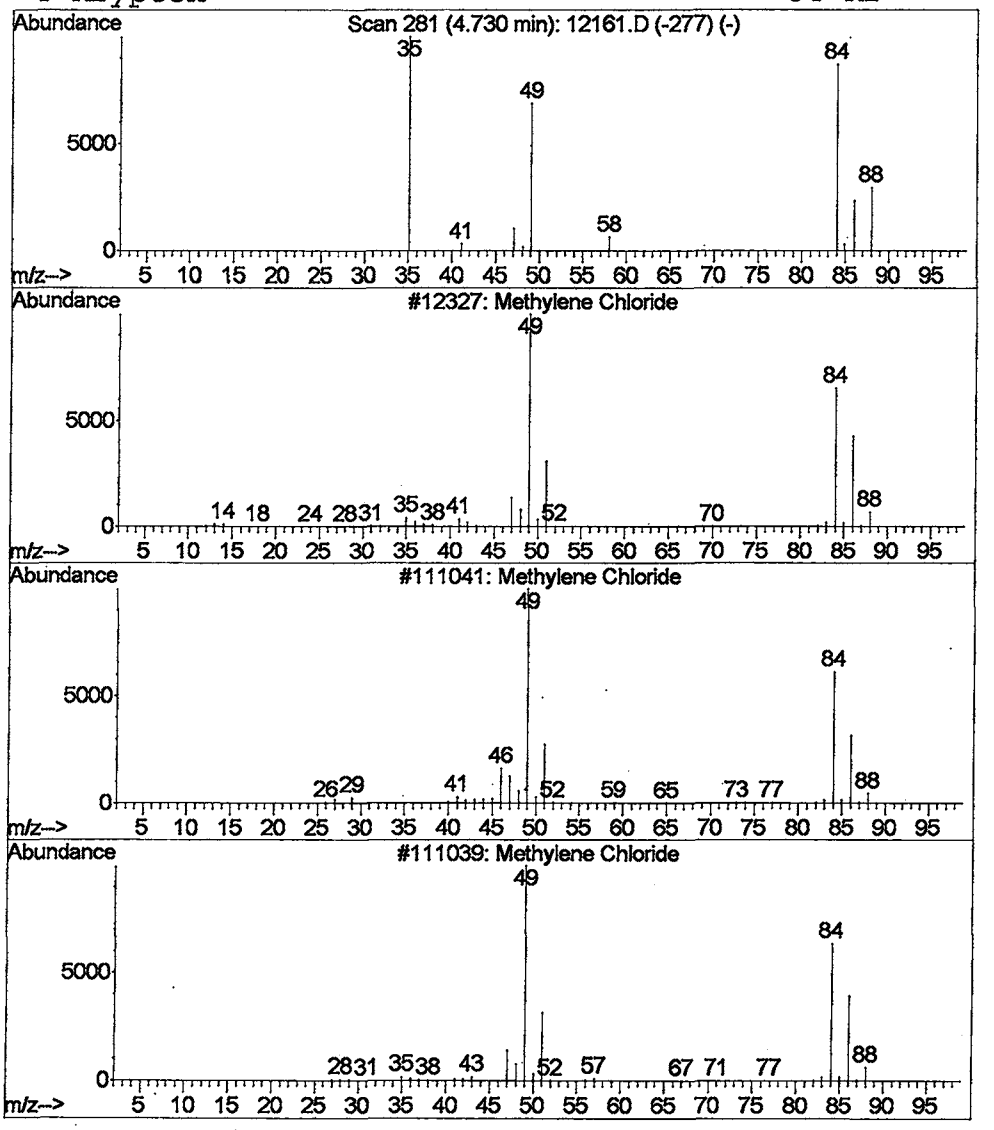
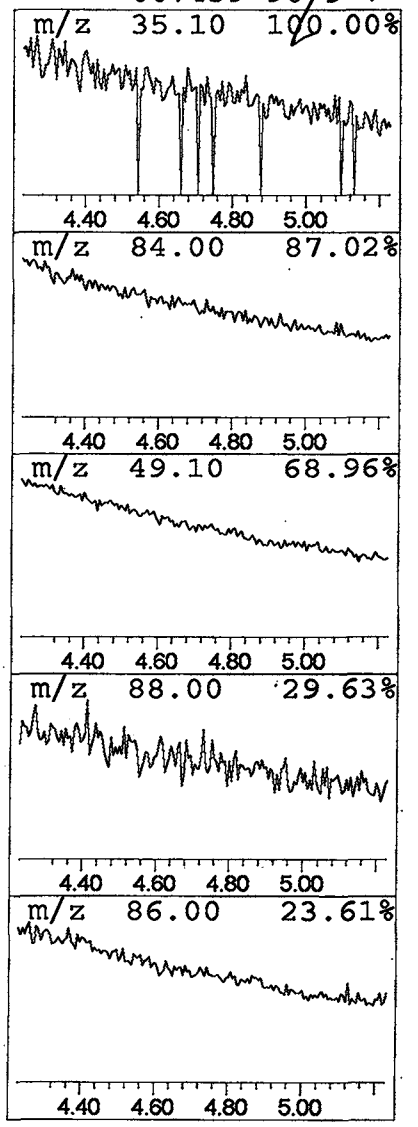
Peak Number 7 Methylene Chloride Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.73	0.60 ug/L	336827	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methylene Chloride	84	CH2Cl2	000075-09-2	10
2	Methylene Chloride	84	CH2Cl2	000075-09-2	9
3	Methylene Chloride	84	CH2Cl2	000075-09-2	9
4	Krypton	84	Kr	007439-90-9	7





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\12161.D
Acq On : 23 May 2002 8:33 pm
Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 14
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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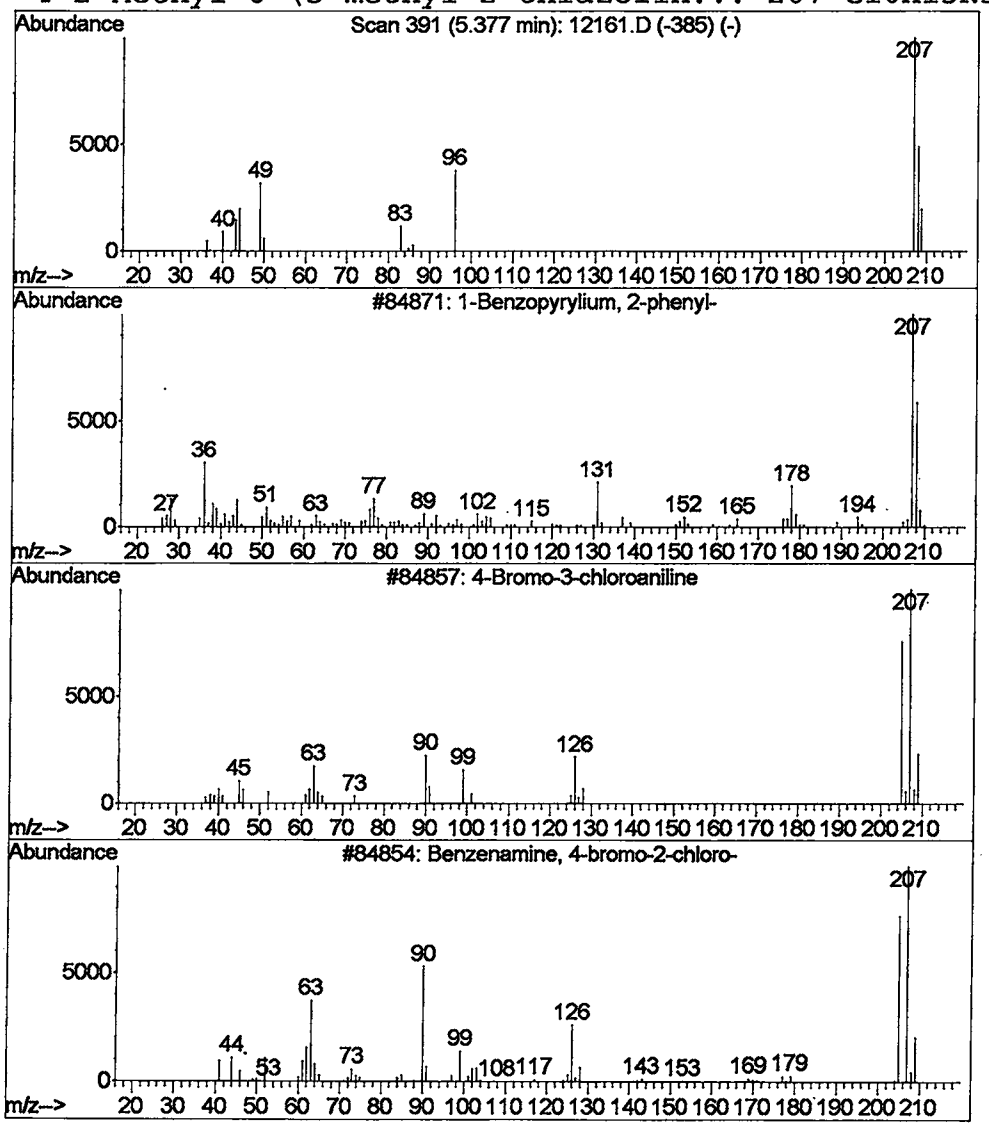
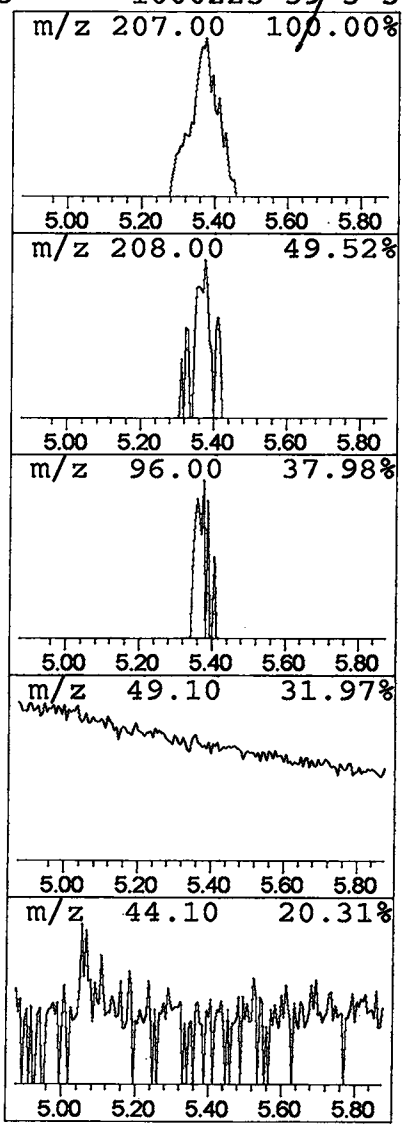
Peak Number 8 1-Benzopyrylium, 2-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.38	0.29 ug/L	160748	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Benzopyrylium, 2-phenyl-	207	C15H11O	014051-53-7	5
2	4-Bromo-3-chloroaniline	205	C6H5BrClN	021402-26-6	5
3	Benzenamine, 4-bromo-2-chloro-	205	C6H5BrClN	038762-41-3	5
4	2-Methyl-6-(5-methyl-2-thiazolin...	207	C10H13N3S	1000225-39-3	5





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\12161.D
Acq On : 23 May 2002 8:33 pm
Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 14
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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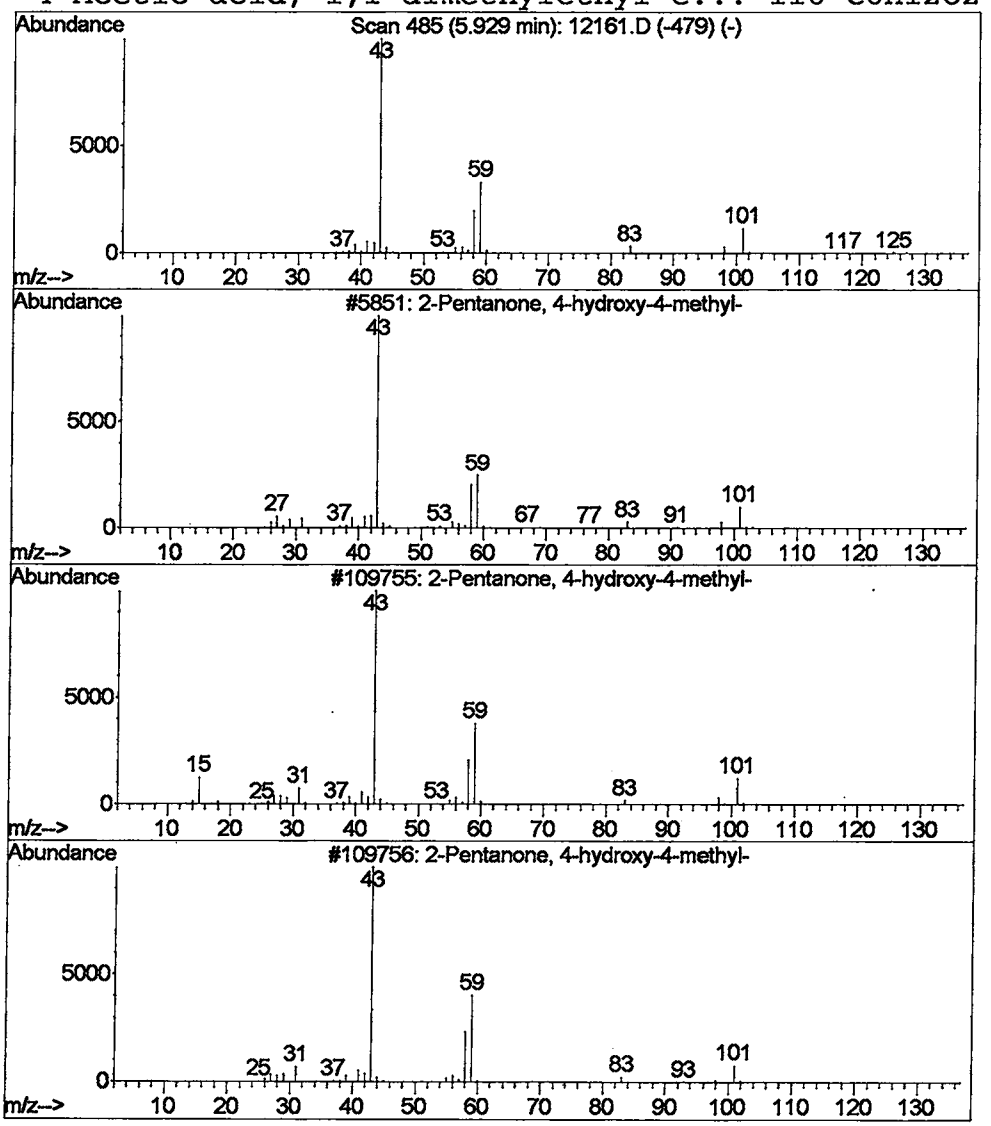
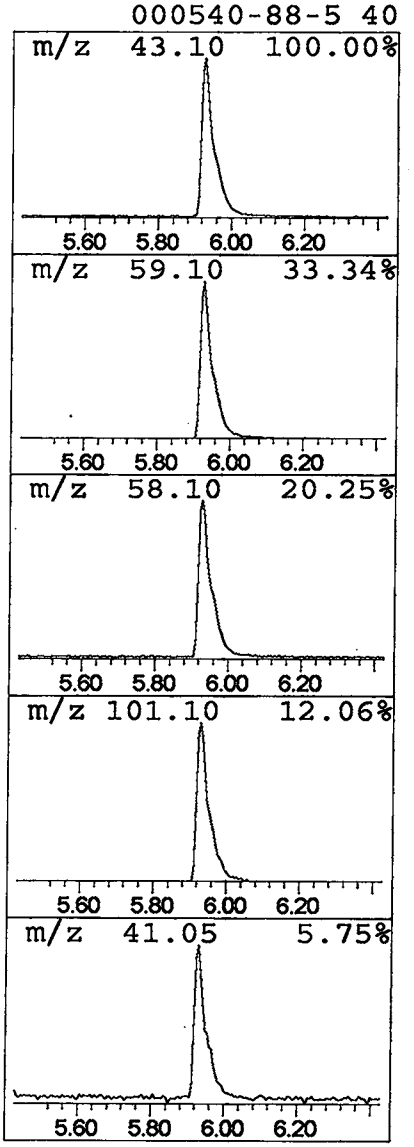
Peak Number 9 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.93	9.72 ug/L	5455810	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	40





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\12161.D
Acq On : 23 May 2002 8:33 pm
Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 14
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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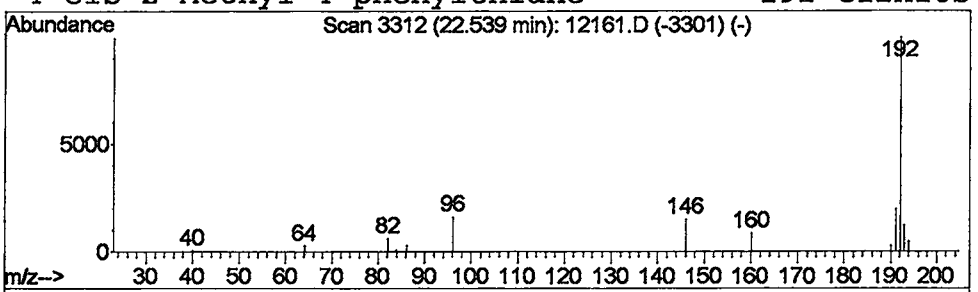
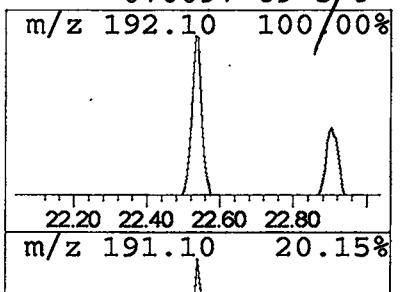
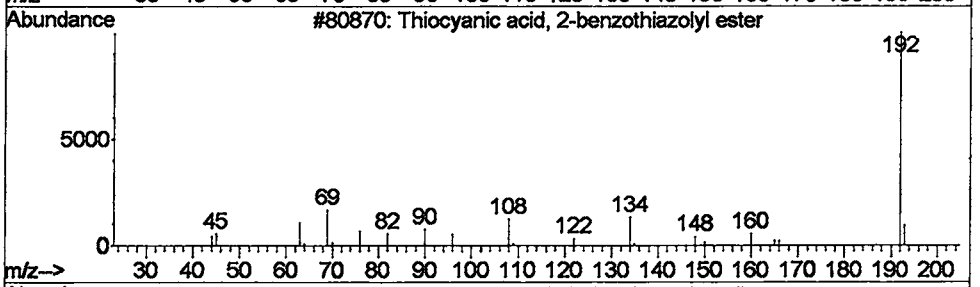
Peak Number 10 Thiocyanic acid, 2-benzothi... Concentration Rank 11

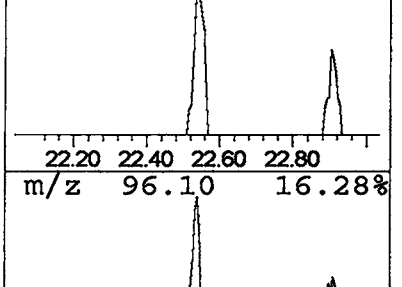
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.54	0.25 ug/L	207181	Phenanthranene-d10	22.91

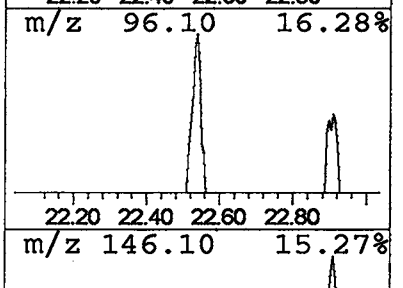
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thiocyanic acid, 2-benzothiazoly...	192	C8H4N2S2	006011-99-0	42
2	2-Propenal, 3-(3,4-dimethoxyphen...	192	C11H12O3	004497-40-9	39
3	1,5-Naphthalenedithiole	192	C10H8S2	1000223-69-5	39
4	cis-2-Methyl-4-phenylthiane	192	C12H16S	076097-69-3	9

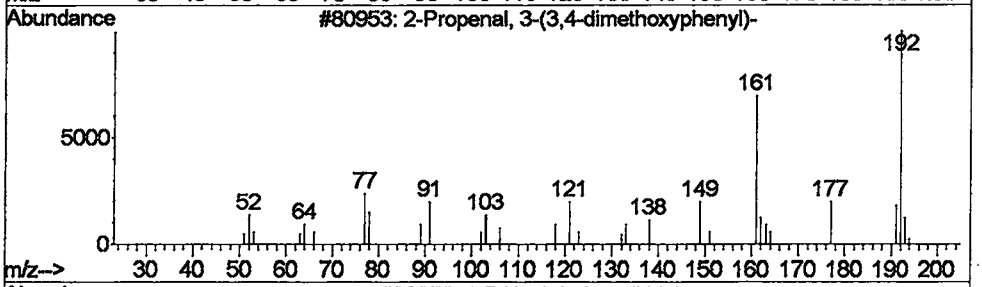


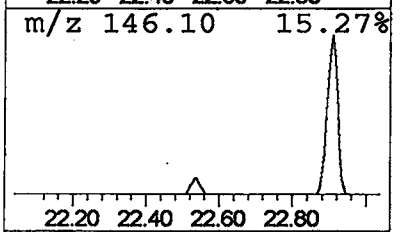


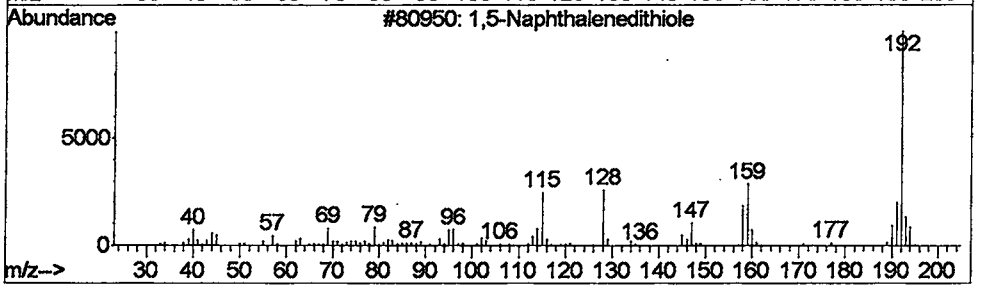


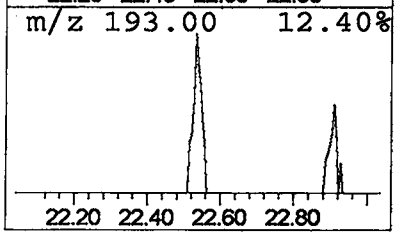












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\12161.D Vial: 14
Acq On : 23 May 2002 8:33 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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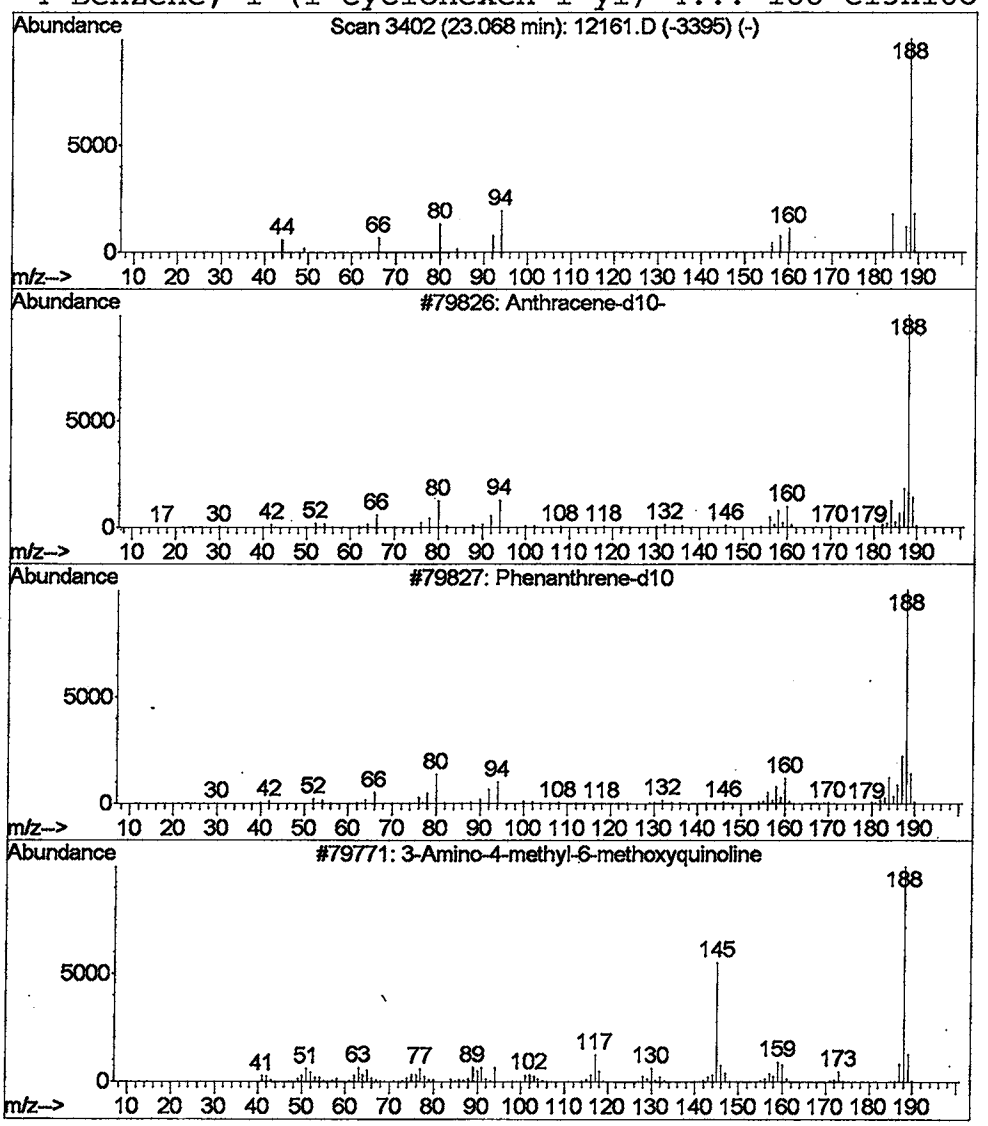
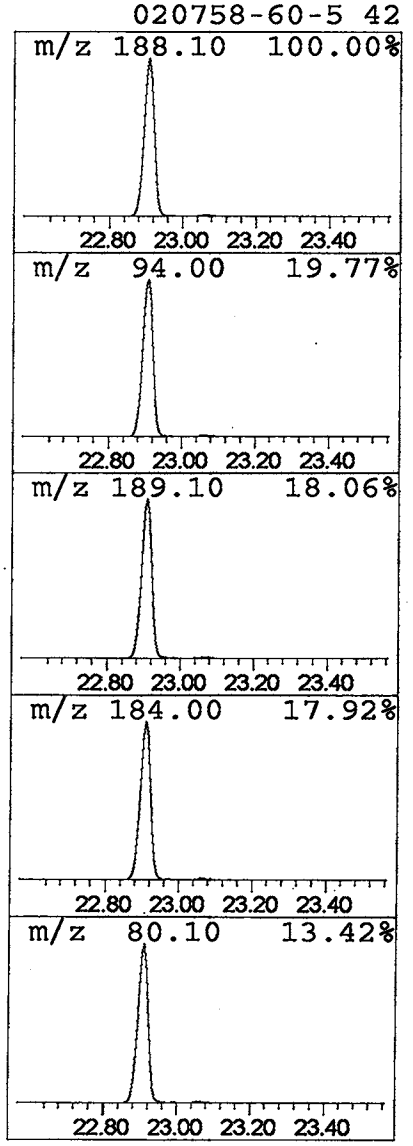
Peak Number 11 Anthracene-d10- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.07	0.26 ug/L	223480	Phenanthranene-d10	22.91

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-d10-	188	C14D10	001719-06-8	83
2	Phenanthrene-d10	188	C14D10	001517-22-2	80
3	3-Amino-4-methyl-6-methoxyquinoline	188	C11H12N2O	1000213-71-3	42
4	Benzene, 1-(1-cyclohexen-1-yl)-4...	188	C13H16O	020758-60-5	42





Tentatively Identified Compound (LSC) summary

Operator ID: Mclean Date Acquired: 23 May 2002 8:33 pm
Data File: C:\MSDCHEM\1\DATA\052302\12161.D
Name: 5/22ad
Misc:
Method: C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title: 508 Modified GC/MS scan
Library Searched: C:\DATABASE\NIST98.L

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--- #	RT	Resp	Conc
Methylene Chloride	3.75	5.3	ug/L	2958380	1	9.90	22453600	40.0
Krypton	3.81	1.0	ug/L	550022	1	9.90	22453600	40.0
1,3-Dioxol-2-one	3.83	2.6	ug/L	1448450	1	9.90	22453600	40.0
2-Chloroethanol	3.91	1.3	ug/L	730513	1	9.90	22453600	40.0
Thiophene	3.95	0.8	ug/L	444049	1	9.90	22453600	40.0
Ethyl Chloride	3.98	3.0	ug/L	1696310	1	9.90	22453600	40.0
Methylene Chloride	4.73	0.6	ug/L	336827	1	9.90	22453600	40.0
1-Benzopyrylium, ...	5.38	0.3	ug/L	160748	1	9.90	22453600	40.0
2-Pentanone, 4-hy...	5.93	9.7	ug/L	5455810	1	9.90	22453600	40.0
Thiocyanic acid, ...	22.54	0.2	ug/L	207181	4	22.91	33748100	40.0
Anthracene-d10-	23.07	0.3	ug/L	223480	4	22.91	33748100	40.0