User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 06/04/2002 Time: 15:05:19

Sample: AE11927
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 6/4/02 15:04 Ended: 6/4/02 15:04 Analyst: DLV
Page: 1

	Component Name	Vol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		102		10.0

Comments for Sample AE11927 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 06-04-2002 Time: 15:05:44

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\11927.D Vial: 7
Acq On : 23 May 2002 2:51 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 29 09:44:48 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	4197202	40.00	ug/L	0.00
10) Napthalene-d8	13.39	136	16686282	40.00	ug/L	-0.01
17) Acenapthalene-d10	18.53	164	9189176	40.00	ug/L	0.00
29) Phenanthranene-d10	22.91	188	14304496	40.00	ug/L	0.00
37) Chrysene-d12	30.76	240	9441484	40.00	ug/L	-0.04
43) Perylene-d12	34.66	264	4949641	40.00	ug/L	-0.02
System Monitoring Compounds						
27) TCMX	20.51	209	1731521	40.83	ug/L	0.00
Spiked Amount 40.000			Recovery = 102.08%			
Target Compounds						Qvalue
2) n-nitros-dimethyl amine	0.00	74	0	N.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	20.10	149	24181	N.D.		
25) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.		
26) Fluorene	0.00	166	0	N.D.		
28) Azobenzene	20.51	77	32127	N.D.		
30) Nnitrosodiphenylamine	20.51	169	32895	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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	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	30.77	228	23665	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

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Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
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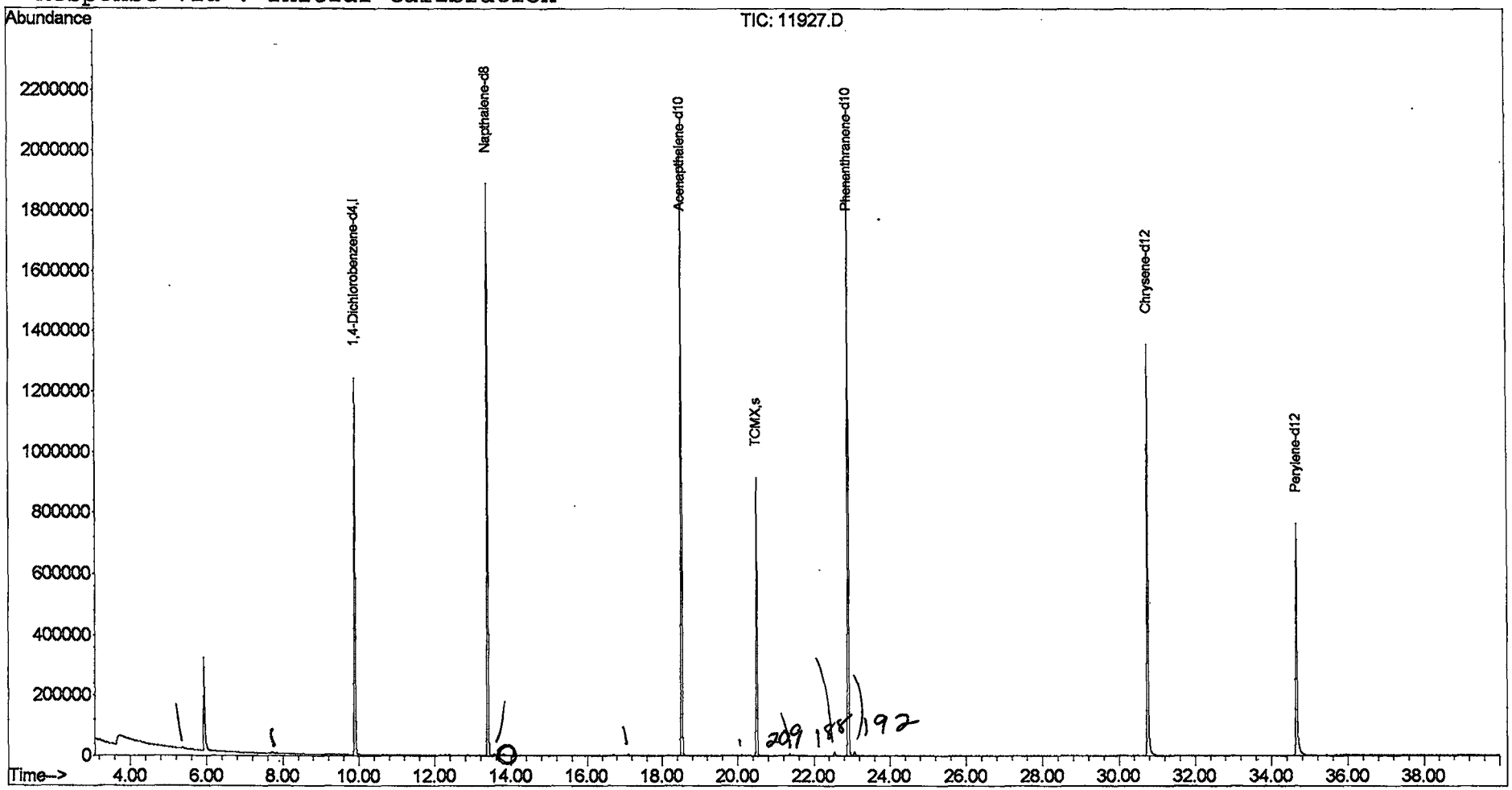


LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\052302\11927.D Vial: 7
Acq On : 23 May 2002 2:51 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

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

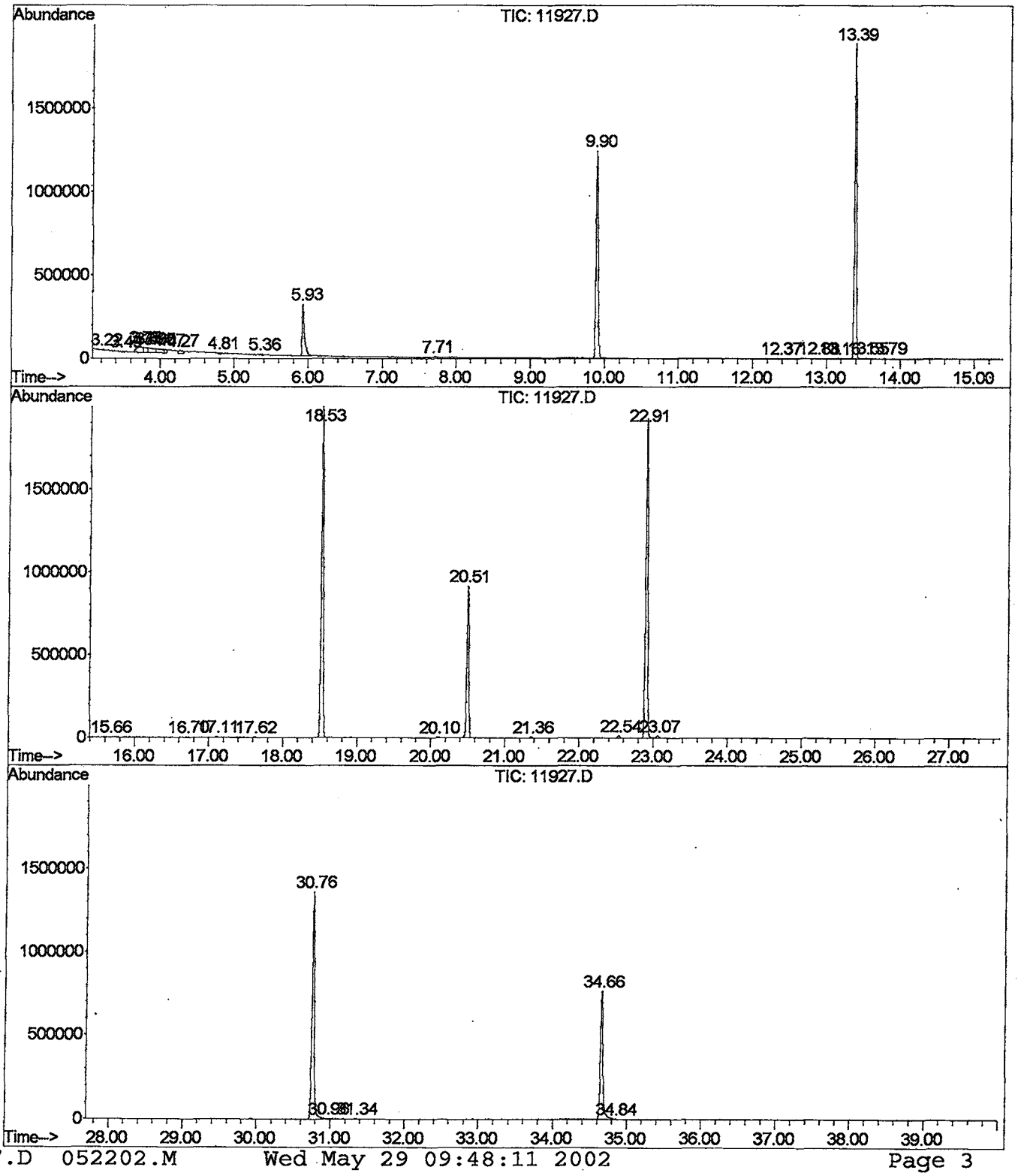
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.226	22	25	30	PV 6	2890	41730	0.11%	0.019%
2	3.490	69	70	75	PV 5	2038	24617	0.06%	0.011%
3	3.749	91	114	117	PV 5	31218	1855305	4.79%	0.864%
4	3.778	117	119	127	VV 5	29522	986455	2.55%	0.459%
5	3.837	127	129	146	VV 7	27672	1681426	4.34%	0.783%
6	3.949	146	148	166	VV 7	23536	1481903	3.83%	0.690%
7	4.072	166	169	172	VV 4	20842	405681	1.05%	0.189%
8	4.266	198	202	211	VV 8	16944	727980	1.88%	0.339%
9	4.812	291	295	297	VV 4	6566	124824	0.32%	0.058%
10	5.365	385	389	398	VV 6	4549	144651	0.37%	0.067%
11	5.929	478	485	513	PV	303158	6601648	17.05%	3.073%
12	7.715	780	789	798	PV 6	4127	82715	0.21%	0.039%
13	9.901	1152	1161	1187	PV	1225377	24959110	64.47%	11.619%
14	12.374	1571	1582	1587	BV 3	1689	18715	0.05%	0.009%
15	12.880	1658	1668	1672	PV 7	1754	49432	0.13%	0.023%
16	13.162	1709	1716	1723	VV 2	2752	43555	0.11%	0.020%
17	13.391	1745	1755	1772	PV	1854279	33918182	87.61%	15.789%
18	13.544	1772	1781	1790	VV 2	5178	106710	0.28%	0.050%
19	13.790	1817	1823	1827	PV 4	4247	58146	0.15%	0.027%
20	15.659	2132	2141	2152	PV 6	1842	50893	0.13%	0.024%
21	16.705	2311	2319	2324	PV 2	2434	37929	0.10%	0.018%
22	17.110	2375	2388	2395	PV 3	5023	75514	0.20%	0.035%
23	17.621	2472	2475	2482	VV 4	2140	31600	0.08%	0.015%
24	18.532	2617	2630	2650	PV	1978120	37556546	97.01%	17.483%
25	20.101	2892	2897	2903	PV 4	3395	65177	0.17%	0.030%
26	20.506	2946	2966	2978	PV	921862	17242786	44.54%	8.027%
27	21.358	3100	3111	3118	BV 2	5949	95314	0.25%	0.044%
28	22.539	3305	3312	3321	PV 3	12596	244254	0.63%	0.114%
29	22.915	3356	3376	3394	PV 2	1919231	38714566	100.00%	18.022%
30	23.068	3394	3402	3411	VV 2	12814	248336	0.64%	0.116%
31	30.765	4697	4712	4744	VV 2	1340959	29026898	74.98%	13.512%
32	30.959	4744	4745	4763	VV 7	2622	88112	0.23%	0.041%
33	31.341	4800	4810	4818	PV 5	2064	39839	0.10%	0.019%
34	34.660	5362	5375	5405	BV 2	766361	17964914	46.40%	8.363%
35	34.843	5405	5406	5411	VV 4	1893	24062	0.06%	0.011%

Sum of corrected areas: 214819523

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\052302\11927.D
Operator : Mclean
Acquired : 23 May 2002 2:51 pm using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 5/22ad
Misc Info :
Vial Number: 7
Quant File :052202.RES (Chemstation Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\11927.D
Acq On : 23 May 2002 2:51 pm
Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 7
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

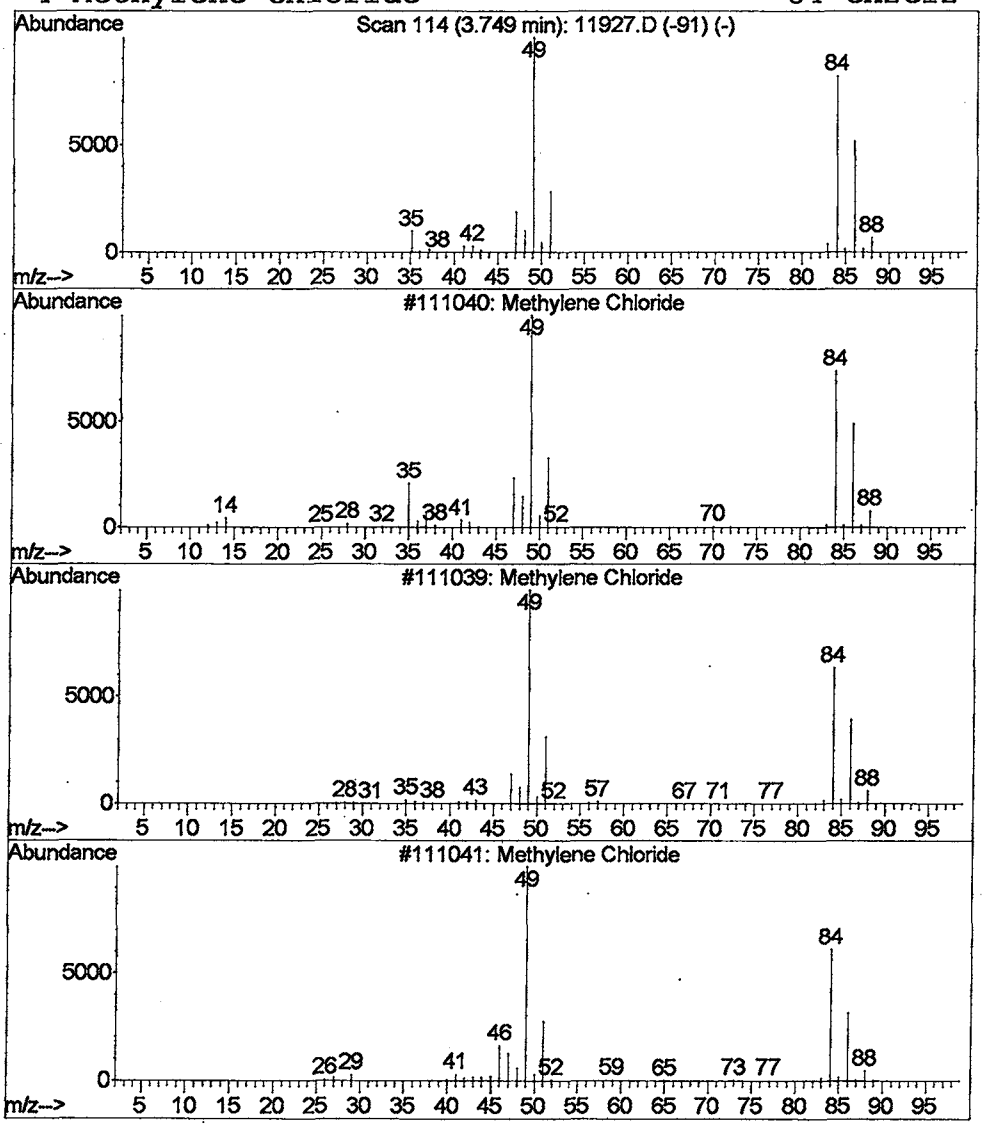
Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

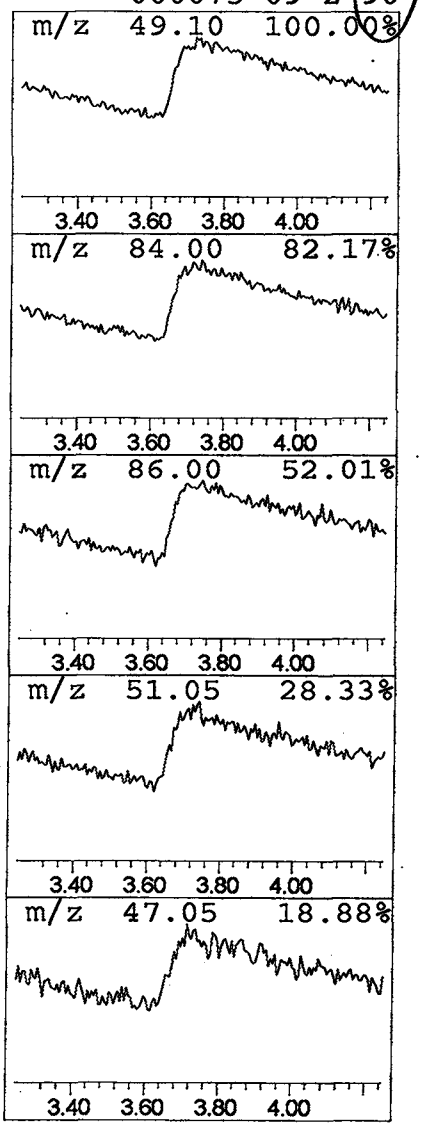
****

Peak Number 1 Methylene Chloride Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.75	2.97 ug/L	1855300	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methylene Chloride	84	CH2Cl2	000075-09-2	94
2	Methylene Chloride	84	CH2Cl2	000075-09-2	91
3	Methylene Chloride	84	CH2Cl2	000075-09-2	90
4	Methylene Chloride	84	CH2Cl2	000075-09-2	90





Library Search Compound Report

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Acq On : 23 May 2002 2:51 pm
Sample : 5/22ad
Misc :
MS Integration Params: LSCINT.e

Vial: 7
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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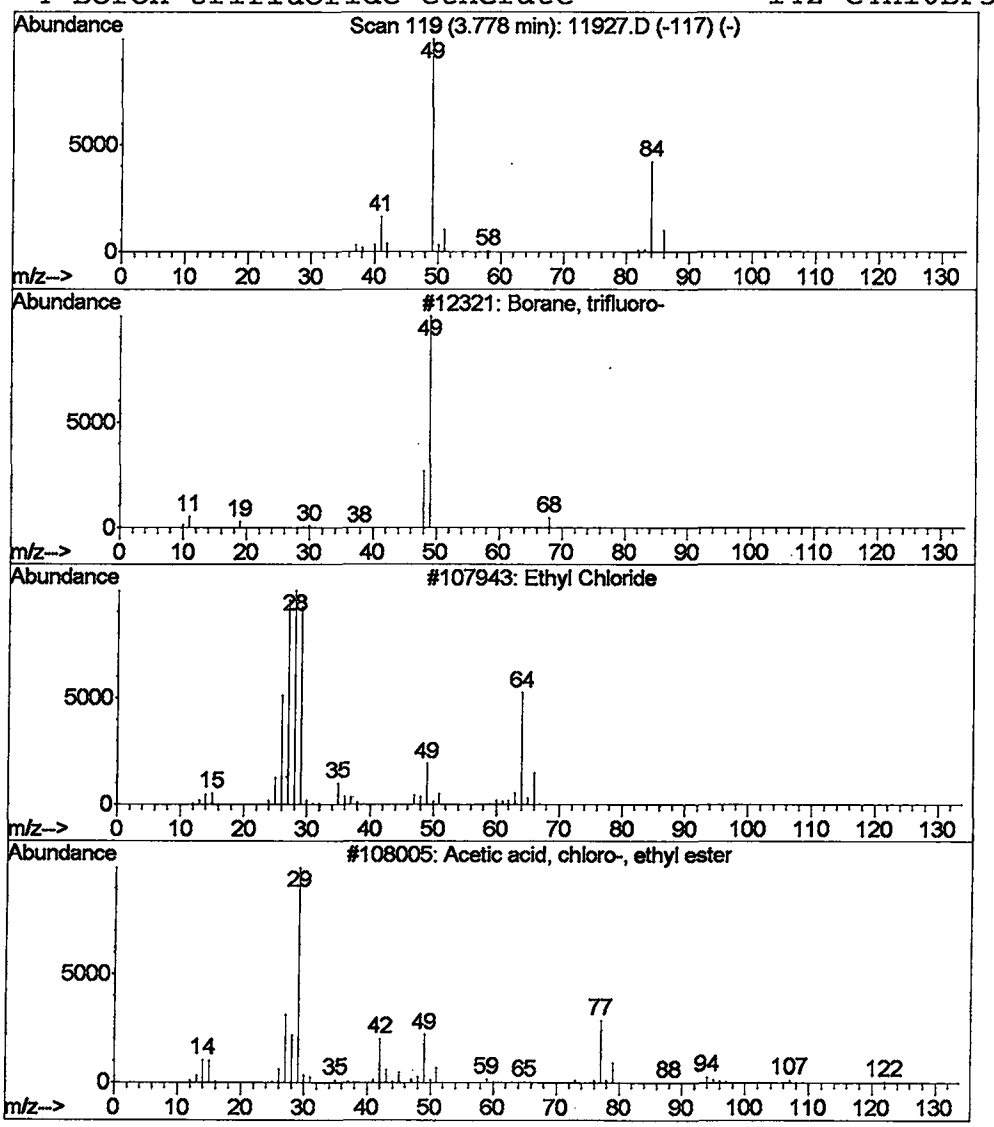
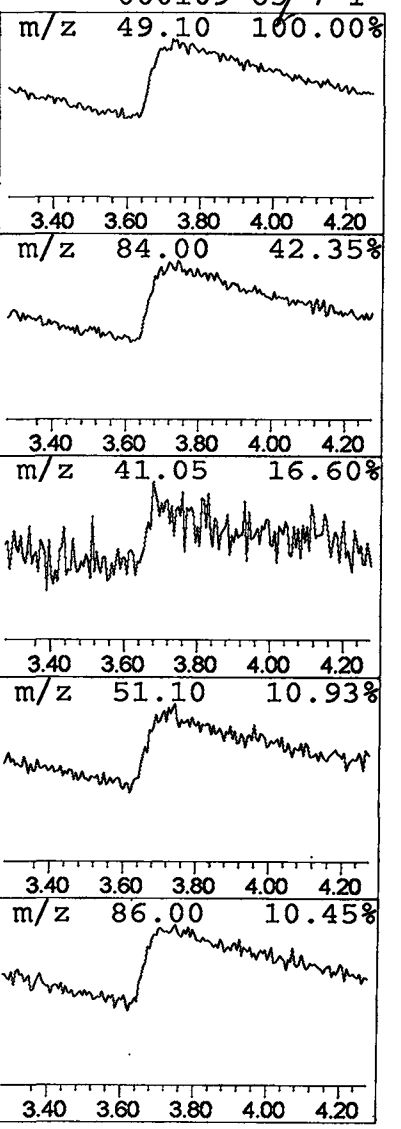
Peak Number 2 Borane, trifluoro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.78	1.58 ug/L	986455	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Borane, trifluoro-	68	BF3	007637-07-2	1
2	Ethyl Chloride	64	C2H5Cl	000075-00-3	1
3	Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	1
4	Boron trifluoride etherate	142	C4H10BF3O	000109-63-7	1





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

Vial: 7
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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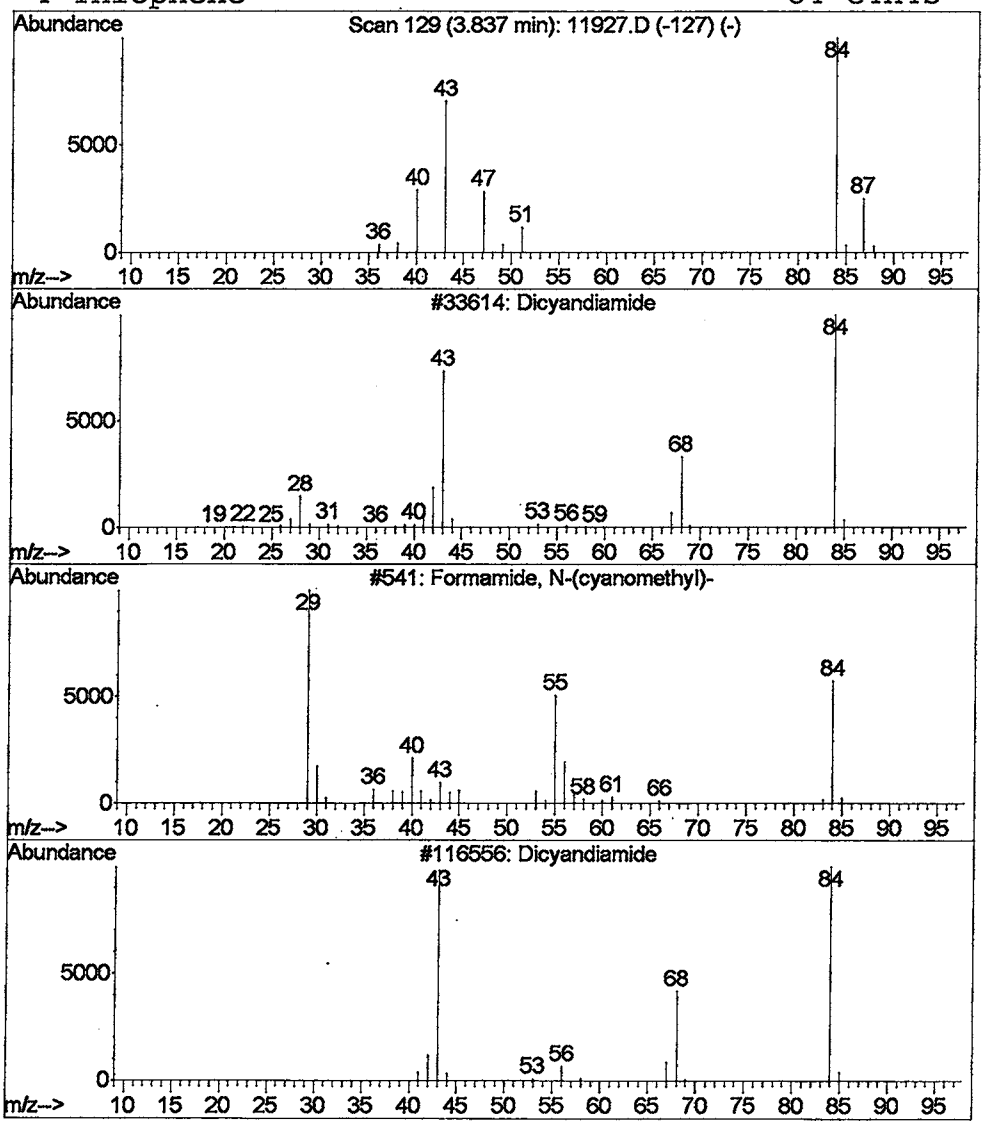
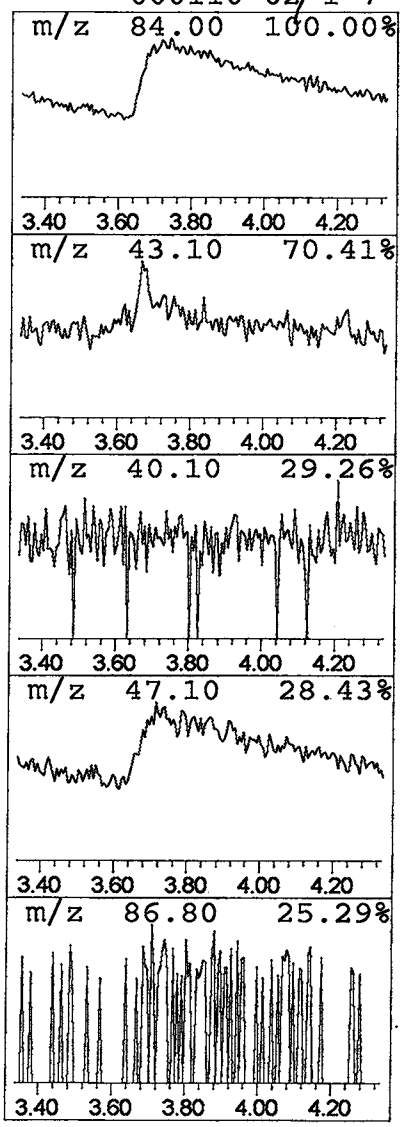
Peak Number 3 Dicyandiamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.83	2.69 ug/L	1681430	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dicyandiamide	84	C2H4N4	000461-58-5	7
2	Formamide, N-(cyanomethyl)-	84	C3H4N2O	005018-27-9	7
3	Dicyandiamide	84	C2H4N4	000461-58-5	7
4	Thiophene	84	C4H4S	000110-02-1	7





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Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
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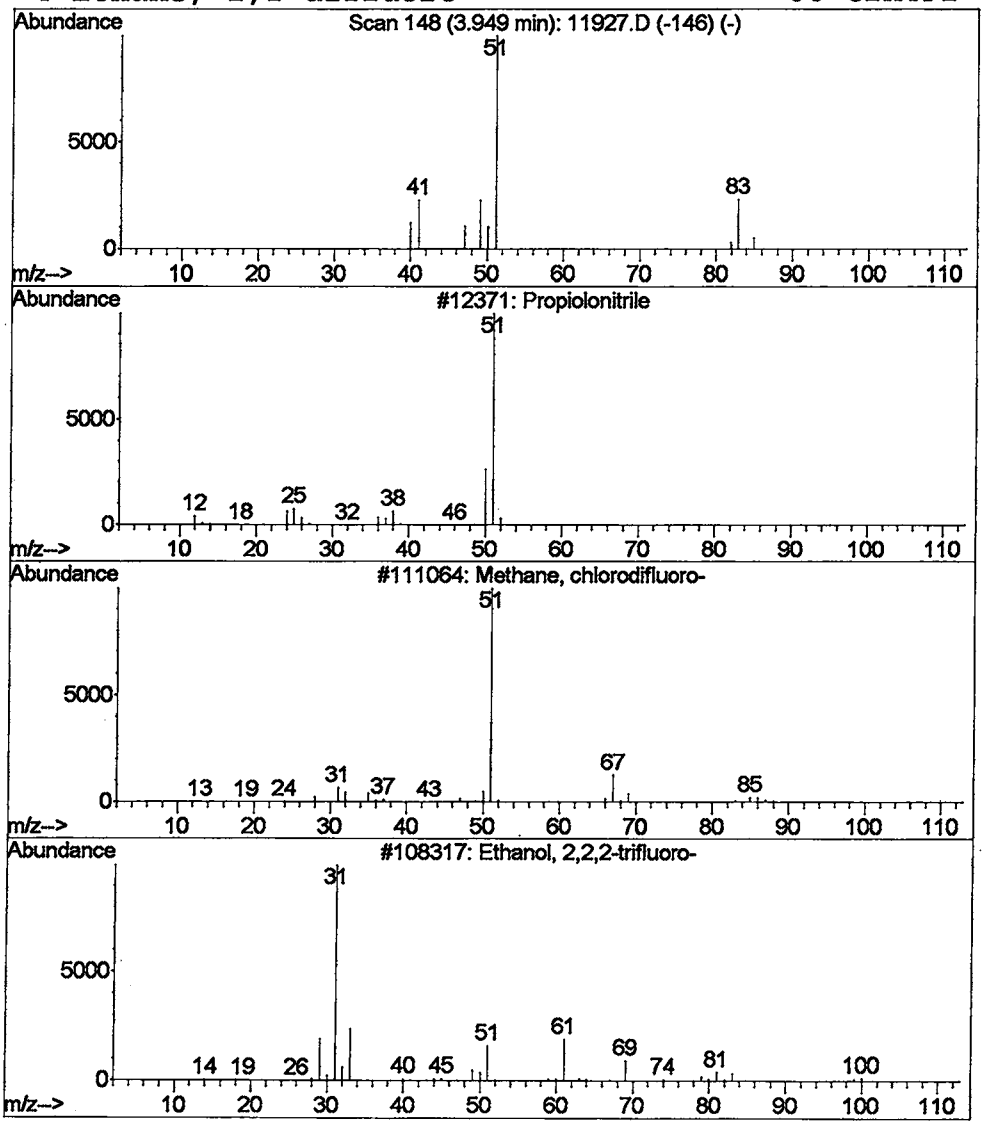
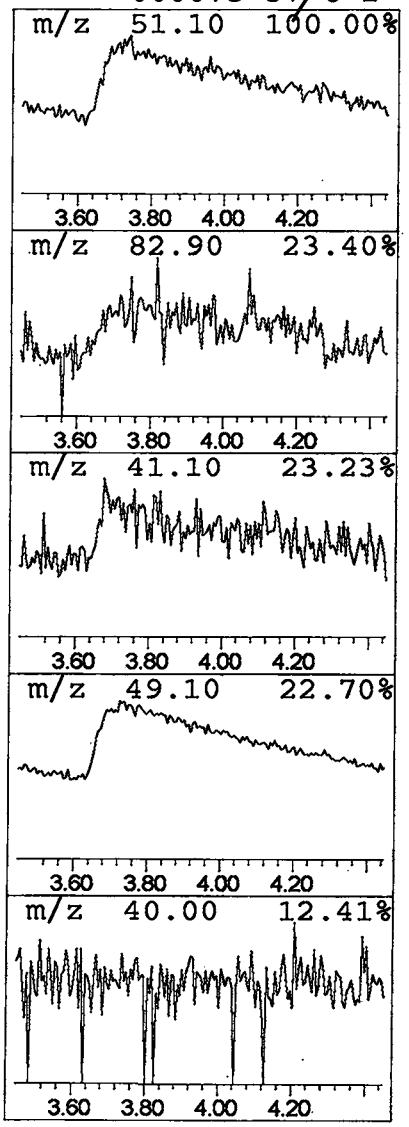
Peak Number 4 Propiolonitrile Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.95	2.37 ug/L	1481900	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propiolonitrile	51	C3HN	001070-71-9	3
2	Methane, chlorodifluoro-	86	CHClF2	000075-45-6	2
3	Ethanol, 2,2,2-trifluoro-	100	C2H3F3O	000075-89-8	2
4	Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	2





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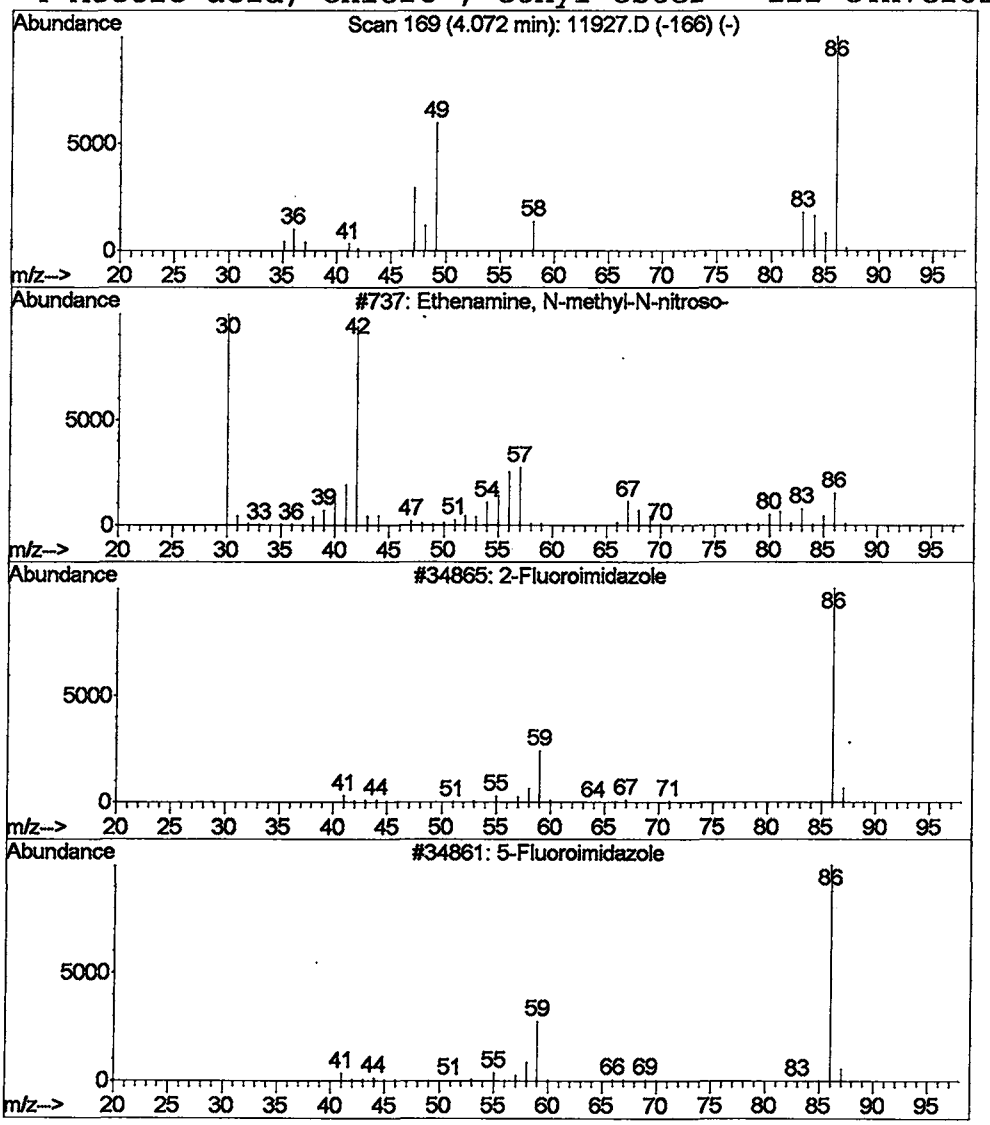
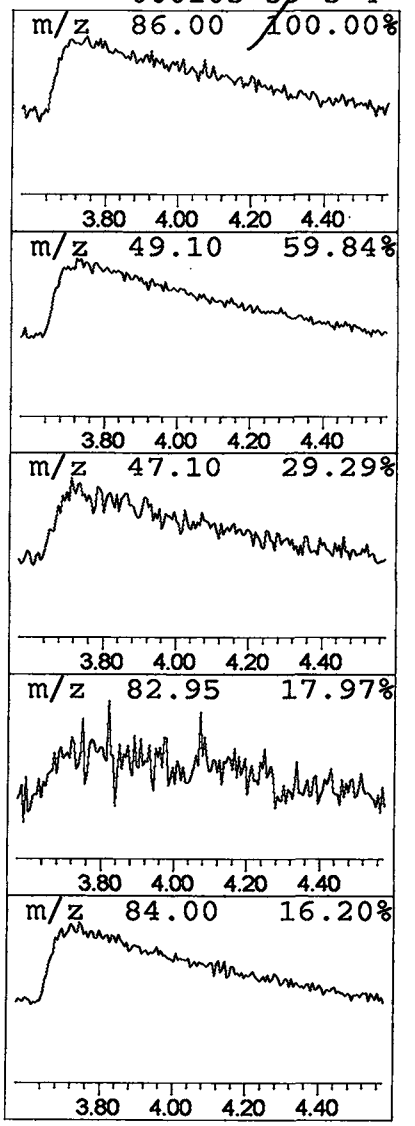
Peak Number 5 Ethenamine, N-methyl-N-nitr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.07	0.65 ug/L	405681	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethenamine, N-methyl-N-nitroso-	86	C3H6N2O	004549-40-0	5
2	2-Fluoroimidazole	86	C3H3FN2	1000129-59-1	5
3	5-Fluoroimidazole	86	C3H3FN2	1000128-06-0	5
4	Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	4





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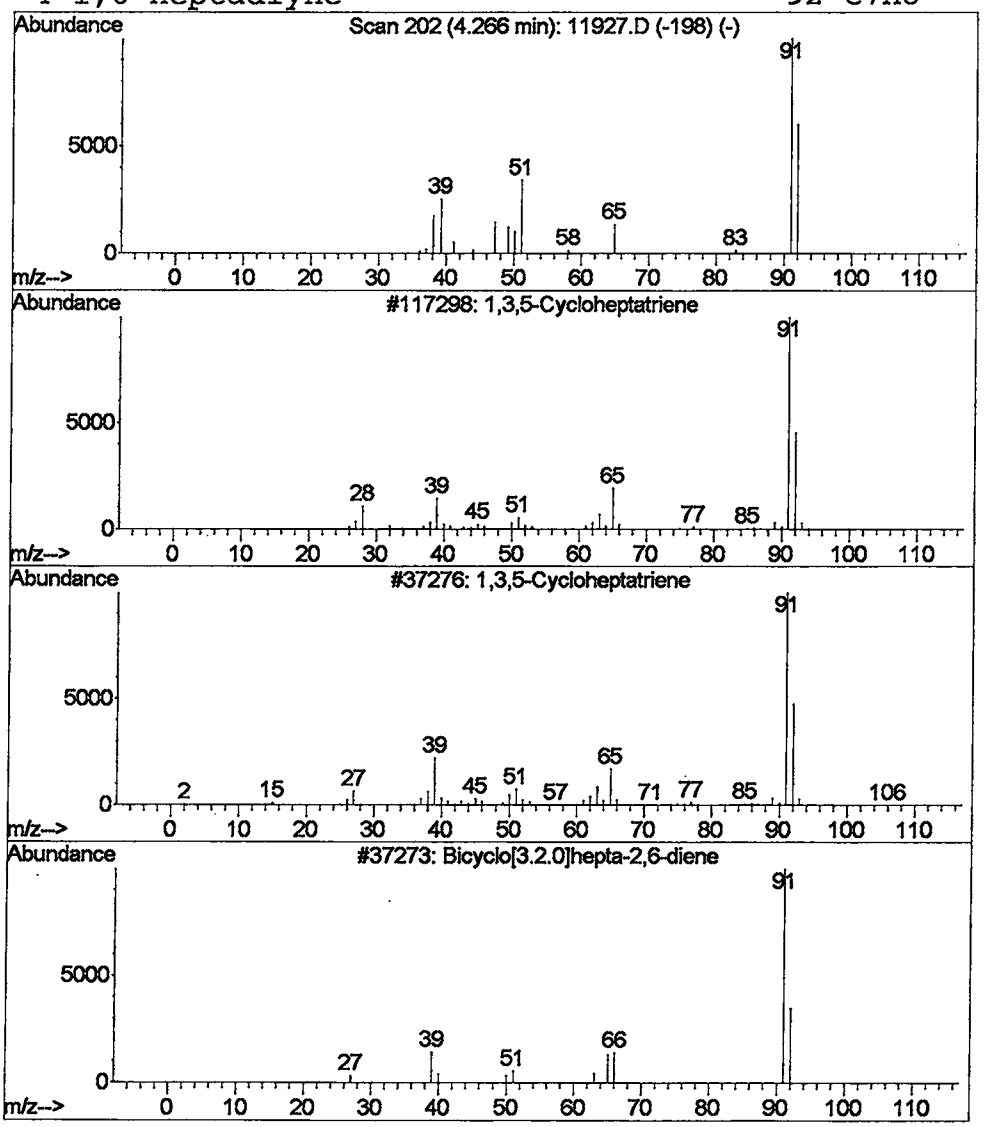
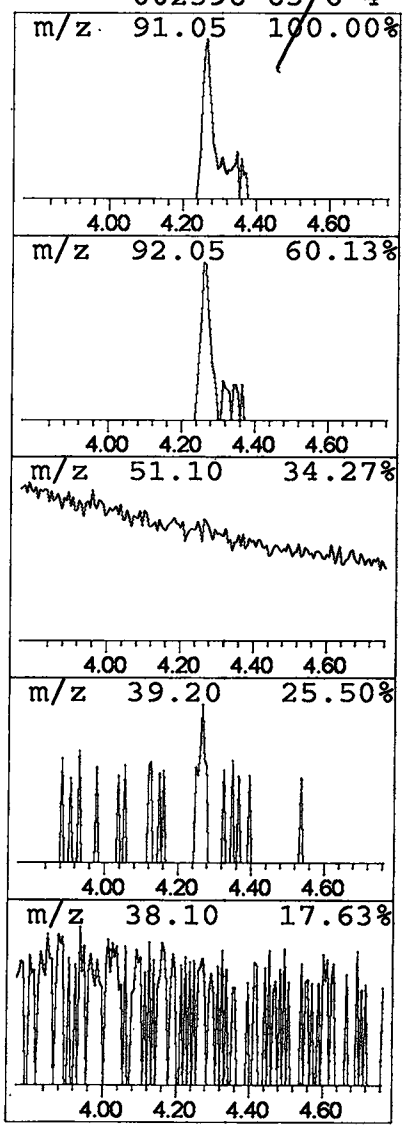
Peak Number 6 1,3,5-Cycloheptatriene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.27	1.17 ug/L	727980	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	7
2	1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	4
3	Bicyclo[3.2.0]hepta-2,6-diene	92	C7H8	002422-86-8	4
4	1,6-Heptadiyne	92	C7H8	002396-63-6	4





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MS Integration Params: LSCINT.e

Vial: 7
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Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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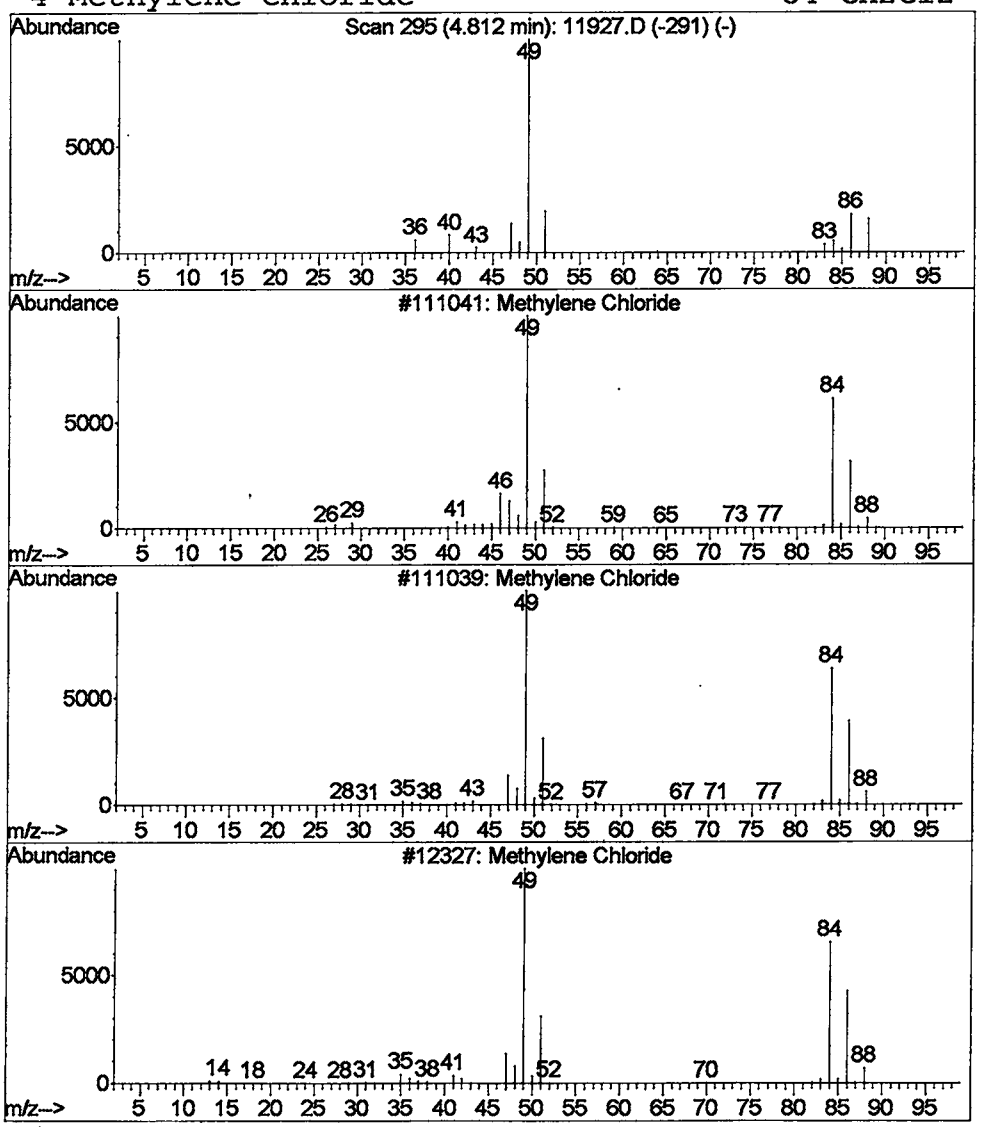
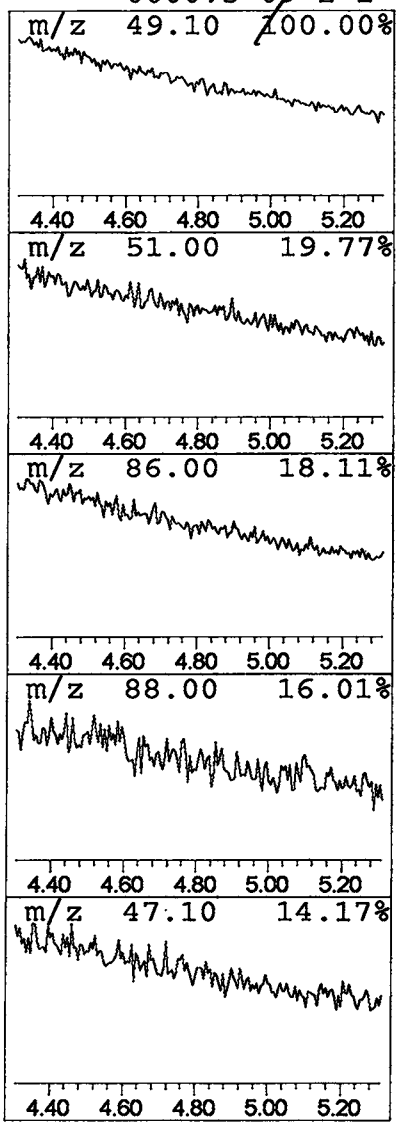
Peak Number 7 Methylene Chloride Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	0.20 ug/L	124824	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methylene Chloride	84	CH2Cl2	000075-09-2	25
2	Methylene Chloride	84	CH2Cl2	000075-09-2	4
3	Methylene Chloride	84	CH2Cl2	000075-09-2	4
4	Methylene Chloride	84	CH2Cl2	000075-09-2	2





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\11927.D Vial: 7
Acq On : 23 May 2002 2:51 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

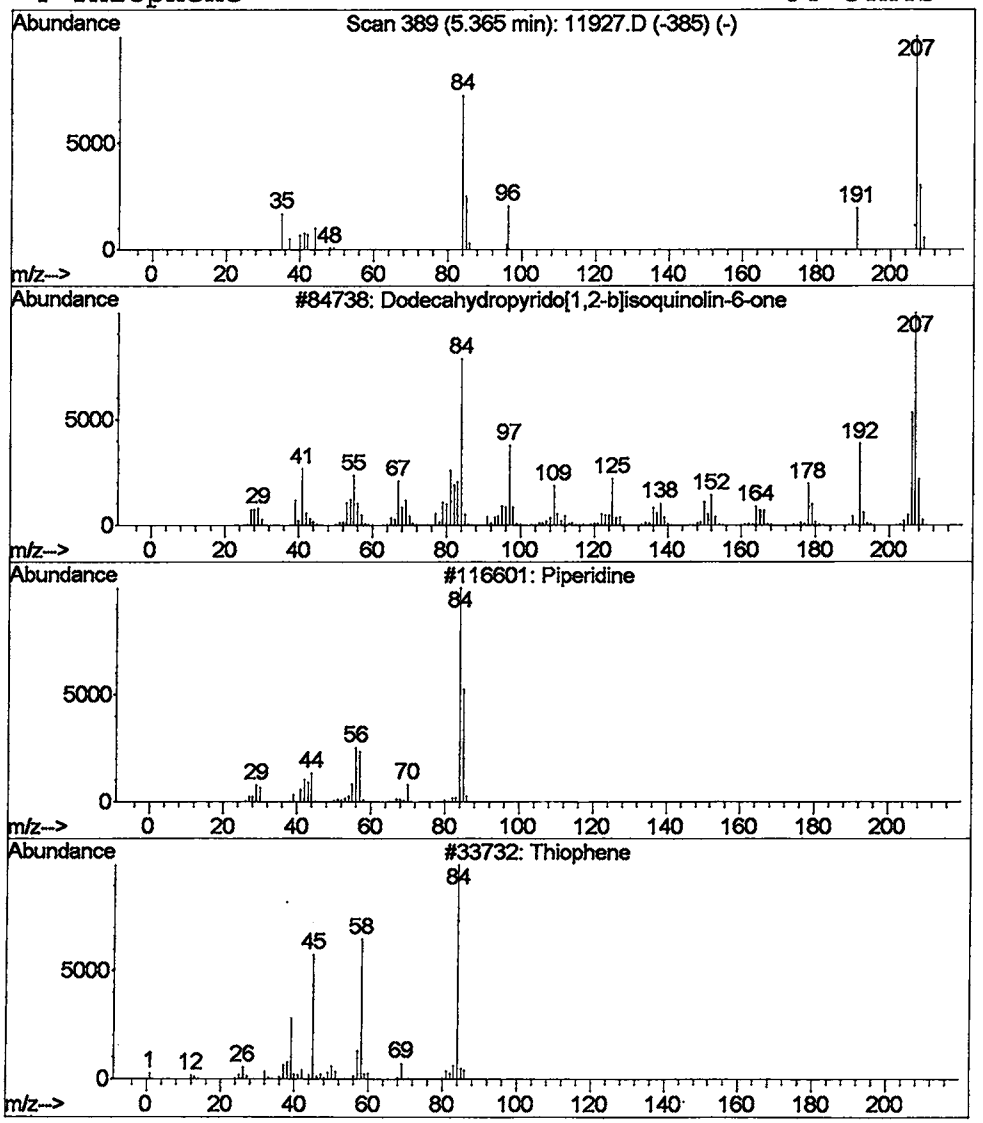
Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

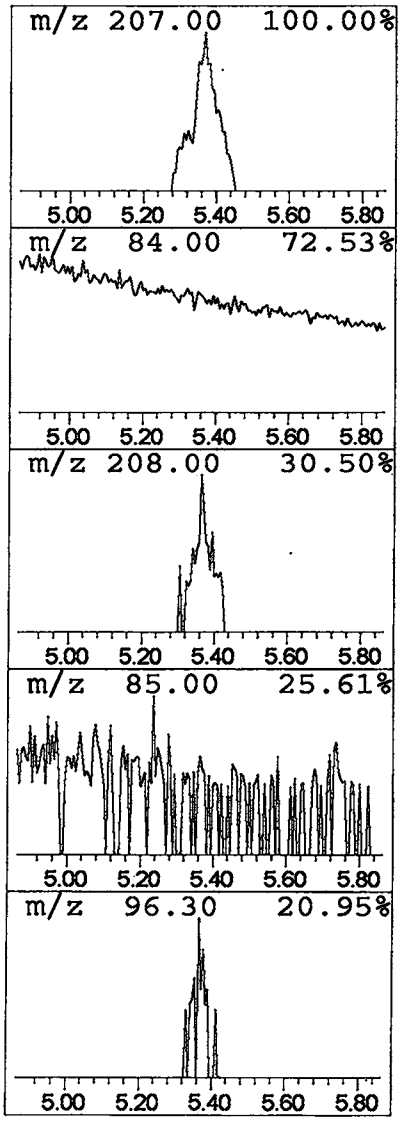
****

Peak Number 8 Dodecahydropyrido[1,2-b]iso... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.36	0.23 ug/L	144651	1,4-Dichlorobenzene-d4	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecahydropyrido[1,2-b]isoquino...	207	C13H21NO	1000195-30-7	38
2	Piperidine	85	C5H11N	000110-89-4	10
3	Thiophene	84	C4H4S	000110-02-1	9
4	Thiophene	84	C4H4S	000110-02-1	9





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\11927.D Vial: 7
Acq On : 23 May 2002 2:51 pm Operator: Mclean
Sample : 5/22ad Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

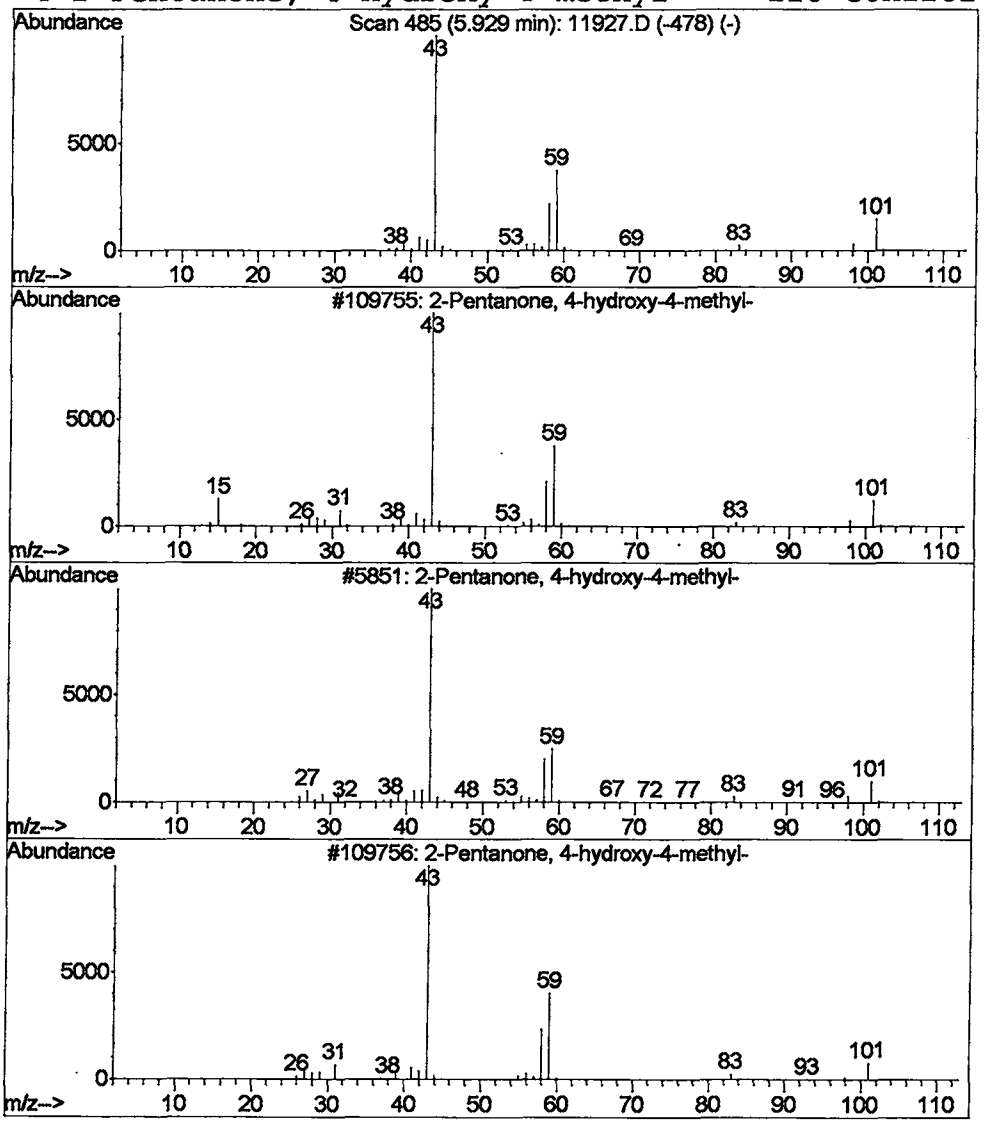
Quant Method : C:\MSDCHEM\1\METHODS\052202.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

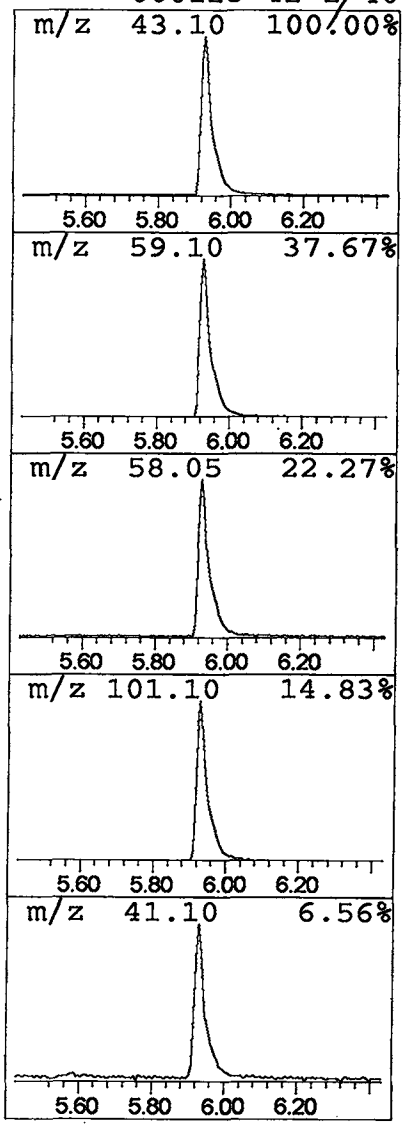
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.93	10.58 ug/L	6601650	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	86
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	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40





Library Search Compound Report

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Acq On : 23 May 2002 2:51 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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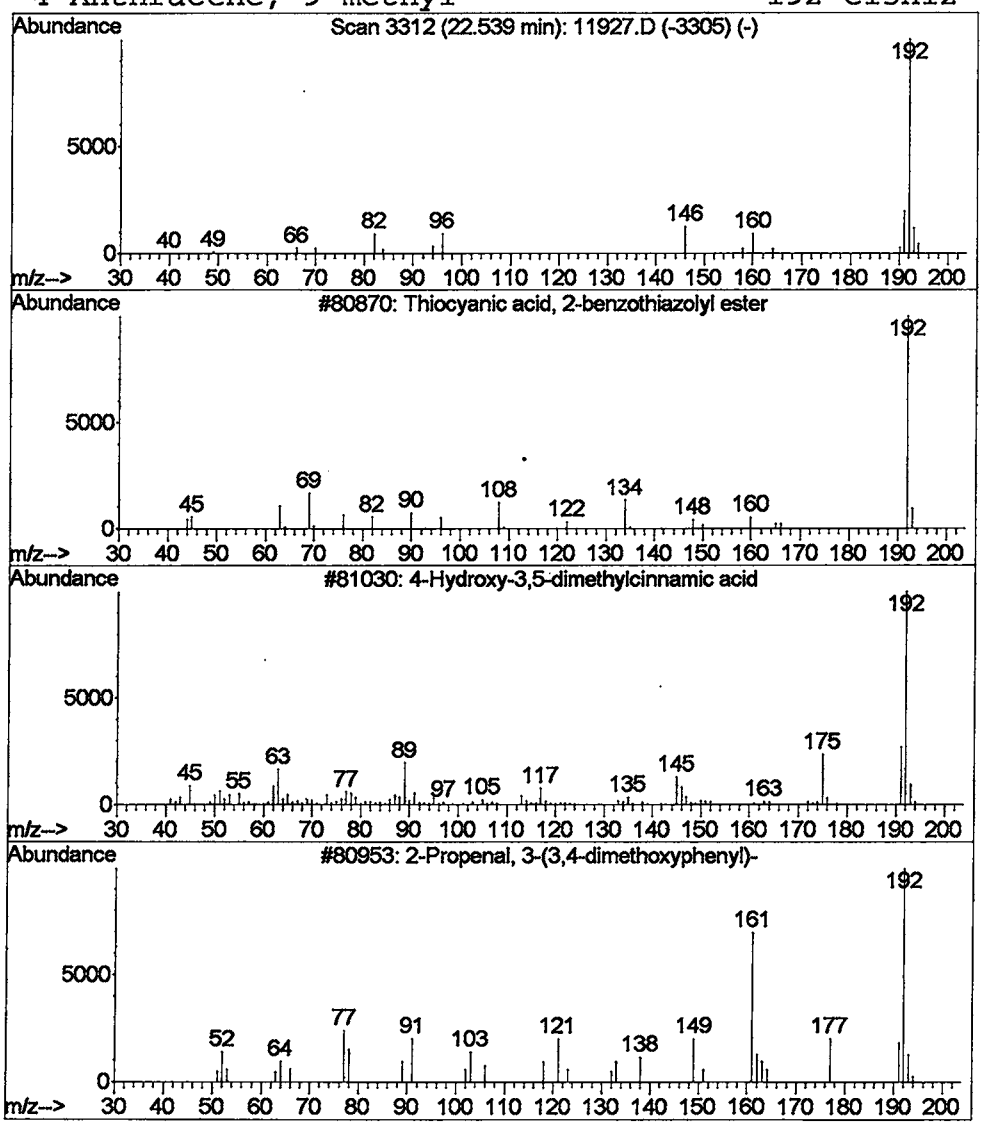
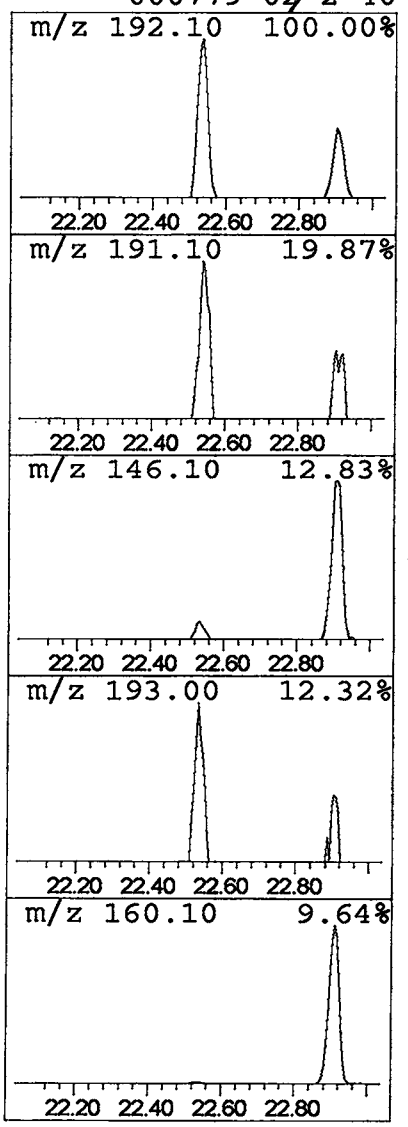
Peak Number 10 Thiocyanic acid, 2-benzothi... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.54	0.25 ug/L	244254	Phenanthranene-d10	22.91

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thiocyanic acid, 2-benzothiazoly...	192	C8H4N2S2	006011-99-0	53
2	4-Hydroxy-3,5-dimethylcinnamic acid	192	C11H12O3	1000132-15-8	45
3	2-Propenal, 3-(3,4-dimethoxyphen...	192	C11H12O3	004497-40-9	42
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	40





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\052302\11927.D Vial: 7
Acq On : 23 May 2002 2:51 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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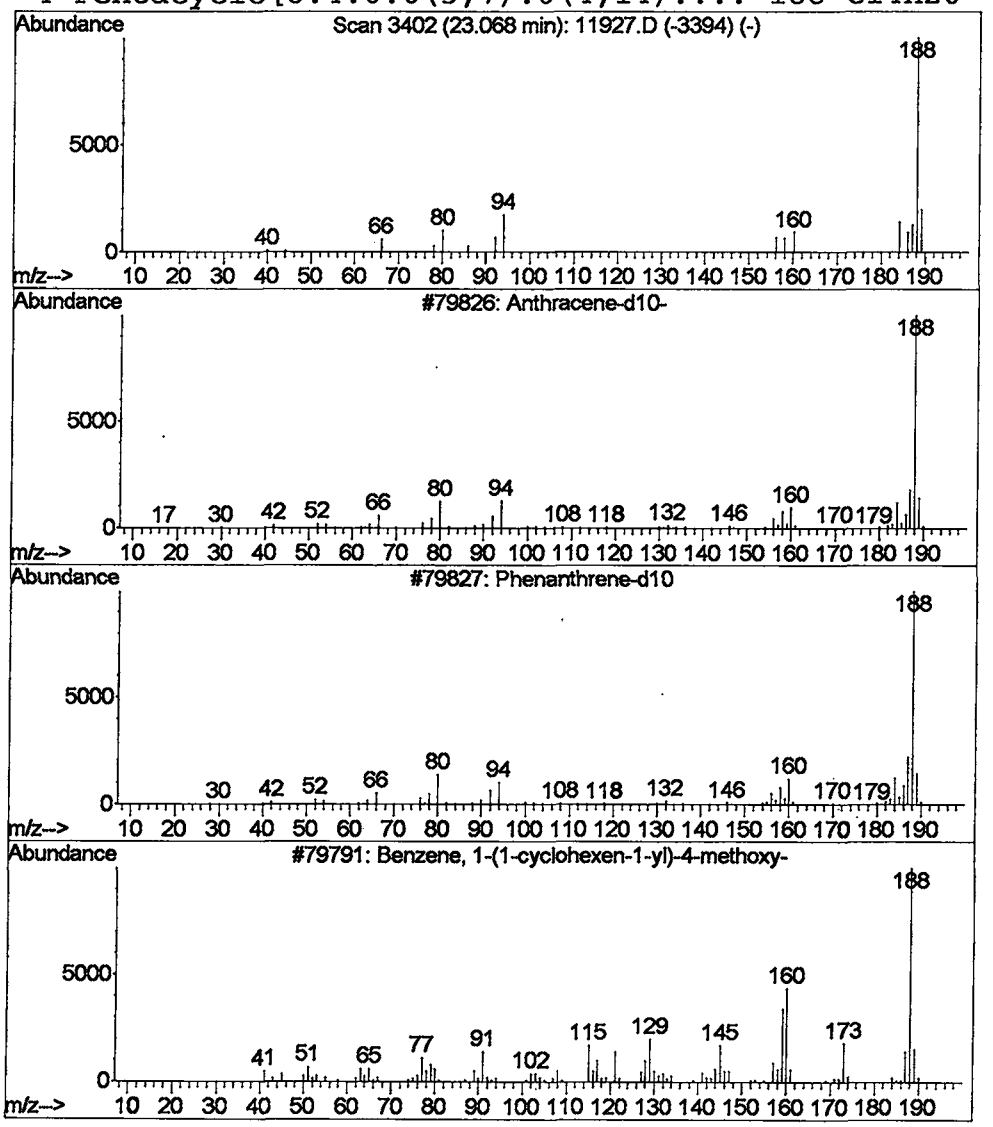
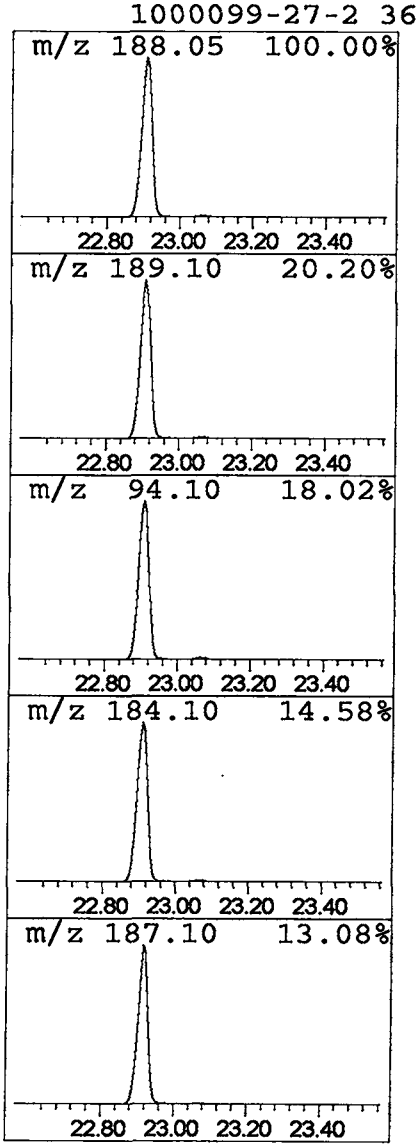
Peak Number 11 Anthracene-d10- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.07	0.26 ug/L	248336	Phenanthranene-d10	22.91

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-d10-	188	C14D10	001719-06-8	91
2	Phenanthrene-d10	188	C14D10	001517-22-2	47
3	Benzene, 1-(1-cyclohexen-1-yl)-4...	188	C13H16O	020758-60-5	42
4	Pentacyclo[8.4.0.0(3,7).0(4,14)....	188	C14H20	1000099-27-2	36





Tentatively Identified Compound (LSC) summary

Operator ID: Mclean Date Acquired: 23 May 2002 2:51 pm
Data File: C:\MSDCHEM\1\DATA\052302\11927.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	3.0	ug/L	1855300	1	9.90	24959100	40.0
Borane, trifluoro-	3.78	1.6	ug/L	986455	1	9.90	24959100	40.0
Dicyandiamide	3.83	2.7	ug/L	1681430	1	9.90	24959100	40.0
Propiolonitrile	3.95	2.4	ug/L	1481900	1	9.90	24959100	40.0
Ethenamine, N-met...	4.07	0.7	ug/L	405681	1	9.90	24959100	40.0
1,3,5-Cycloheptat...	4.27	1.2	ug/L	727980	1	9.90	24959100	40.0
Methylene Chloride	4.81	0.2	ug/L	124824	1	9.90	24959100	40.0
Dodecahydropyrido...	5.36	0.2	ug/L	144651	1	9.90	24959100	40.0
2-Pentanone, 4-hy...	5.93	10.6	ug/L	6601650	1	9.90	24959100	40.0
Thiocyanic acid, ...	22.54	0.3	ug/L	244254	4	22.91	38714600	40.0
Anthracene-d10-	23.07	0.3	ug/L	248336	4	22.91	38714600	40.0