User: Baglioni, Walter
Westchester Dept. of Labs & Research
Date: 02/26/2002 Time: 10:41:28

Sample: AE01136
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 2/25/02 10:41 Ended: 2/26/02 10:41 Analyst: WB
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	n-Nitrosodimethylamine		< MDL		2.0
2	bis(2-Chloroethyl)ether		< MDL		2.0
3	1,3-Dichlorobenzene		< MDL		2.0
4	1,4-Dichlorobenzene		< MDL		2.0
5	1,2-Dichlorobenzene		< MDL		2.0
6	2,2'-oxybis(1-Chloropropane)		< MDL		2.0
7	n-Nitrosodi-n-propylamine		< MDL		2.0
8	Hexachloroethane		< MDL		2.0
9	Nitrobenzene		< MDL		2.0
10	Isophorone		< MDL		2.0
11	bis(2-Chloroethoxy)methane		< MDL		2.0
12	1,2,4-Trichlorobenzene		< MDL		2.0
13	Naphthalene		< MDL		2.0
14	Hexachiorobutadiene		< MDL		2.0
15	2-Chloronaphthalene		< MDL		2.0
16	Dimethylphthalate		< MDL		2.0
17	2,6-Dinitrotoluene		< MDL		2.0
18	Acenaphthylene		< MDL		2.0
19	Acenaphthene		< MDL		2.0
20	2,4-Dinitrotoluene		< MDL		2.0
21	Diethylphthalate		< MDL		2.0
22	4-Chlorophenylphenylether		< MDL		2.0
23	Fluorene		< MDL		2.0
24	Azobenzene/1,2-Diphenylhydra		< MDL		2.0
25	n-Nitrosodiphenylamine		< MDL		2.0
26	4-Bromophenylphenylether		< MDL		2.0
27	Hexachlorobenzene		< MDL		2.0
28	Phenanthrene		< MDL		2.0
29	Anthracene		< MDL		2.0
30	Di-n-butylphthalate		< MDL		2.0
31	Fluoranthene		< MDL		2.0
32	Pyrene		< MDL		2.0
33	Butylbenzylphthalate		< MDL		2.0
34	Benzo(a)anthracene		< MDL		2.0
35	bis(2-Ethylhexyl)phthalate		< MDL		2.0
36	Chrysene		< MDL		2.0
37	Di-n-octylphthalate		< MDL		2.0
38	Benzo(b)fluoranthene		< MDL		2.0
39	Benzo(k)fluoranthene		< MDL		2.0
40	Benzo(a)pyrene		< MDL		2.0
41	Indeno(1,2,3-cd)pyrene		< MDL		2.0
42	Dibenz(a,h)anthracene		< MDL		2.0
43	Benzo(g,h,i)perylene		< MDL		2.0

Data File : C:\MSDCHEM\1\5973N\02FEB25\0116B.D Vial: 1
Acq On : 25 Feb 2002 11:14 am Operator:
Sample : lab blank 1/16 Inst : Instrumen
Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Feb 26 08:18:21 2002 Quant Results File: HP022102.RES

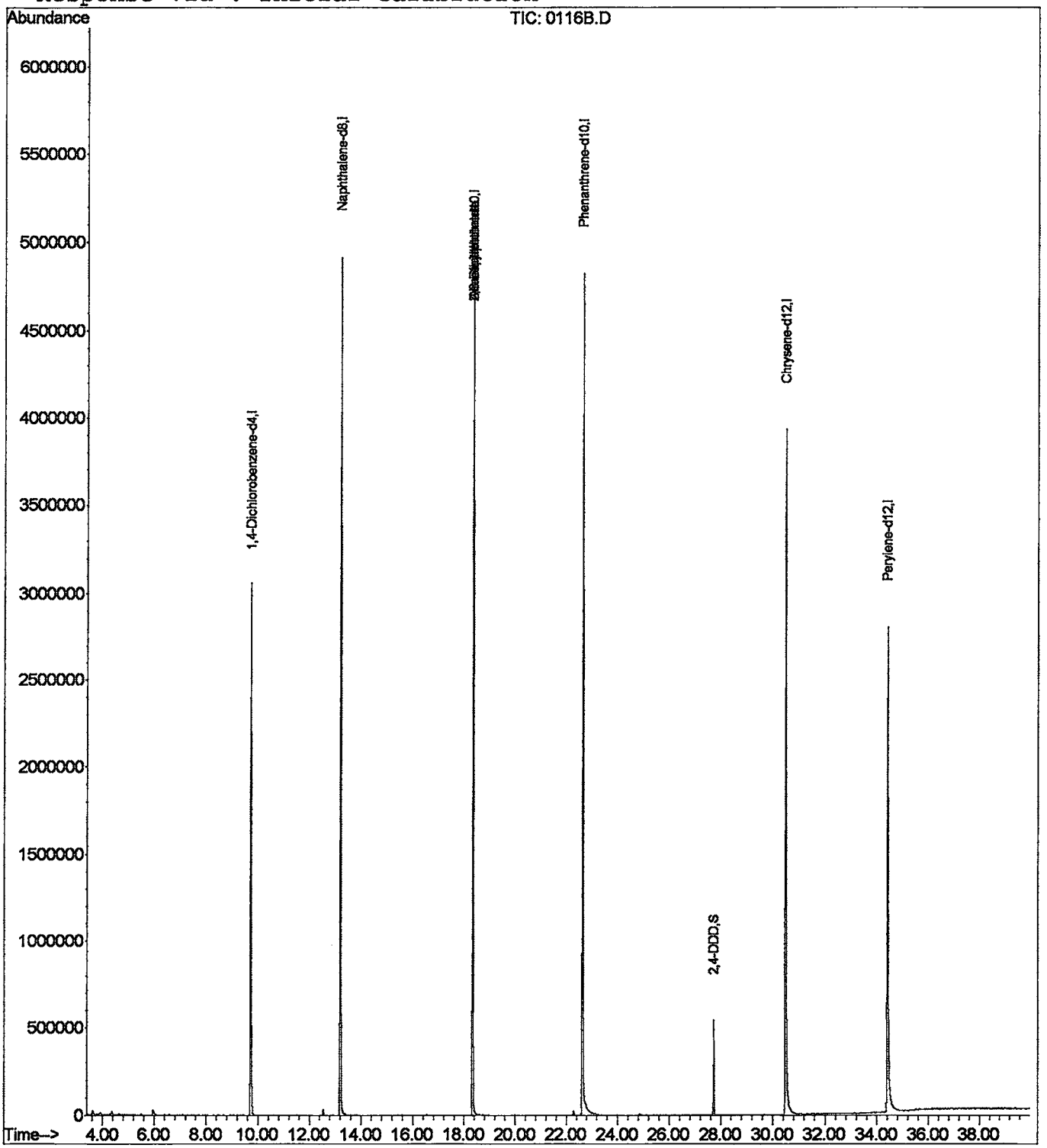
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Last Update : Mon Feb 25 10:47:59 2002
Response via : Initial Calibration
DataAcq Meth : HP020702

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.72	152	1184628	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	5073061	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	2451324	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	4274830	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	3219213	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	2203430	20.00	ug/L	-0.04
System Monitoring Compounds						
37) 2,4-DDD	27.72	235	190006	3.33	ug/L	0.00
Spiked Amount 5.000			Recovery = 66.60%			
Target Compounds						Qvalue
20) Dimethylphthalate	18.33	163	406424	2.73	ug/L	# 58
21) 2,6-Dinitrotoluene	18.33	165	315187	9.18	ug/L	# 27

(#) = qualifier out of range (m) = manual integration
0116B.D HP022102.M Tue Feb 26 08:18:22 2002 Page 1

Data File : C:\MSDCHEM\1\5973N\02FEB25\0116B.D Vial: 1
Acq On : 25 Feb 2002 11:14 am Operator:
Sample : lab blank 1/16 Inst : Instrumen
Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Feb 26 8:18 2002 Quant Results File: HP022102.R

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
Title : 508 scan method
Last Update : Mon Feb 25 10:47:59 2002
Response via : Initial Calibration



Data File : C:\MSDCHEM\1\5973N\02FEB25\0116B.D Vial: 1
Acq On : 25 Feb 2002 11:14 am Operator:
Sample : lab blank 1/16 Inst : Instrumen
Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
Title : 508 scan method
Smoothing : OFF Filtering: 5
Sampling : 2 Min Area: 100 Area counts
Start Thrs: 0.2 Max Peaks: 20
Stop Thrs : 0 Peak Location: TOP

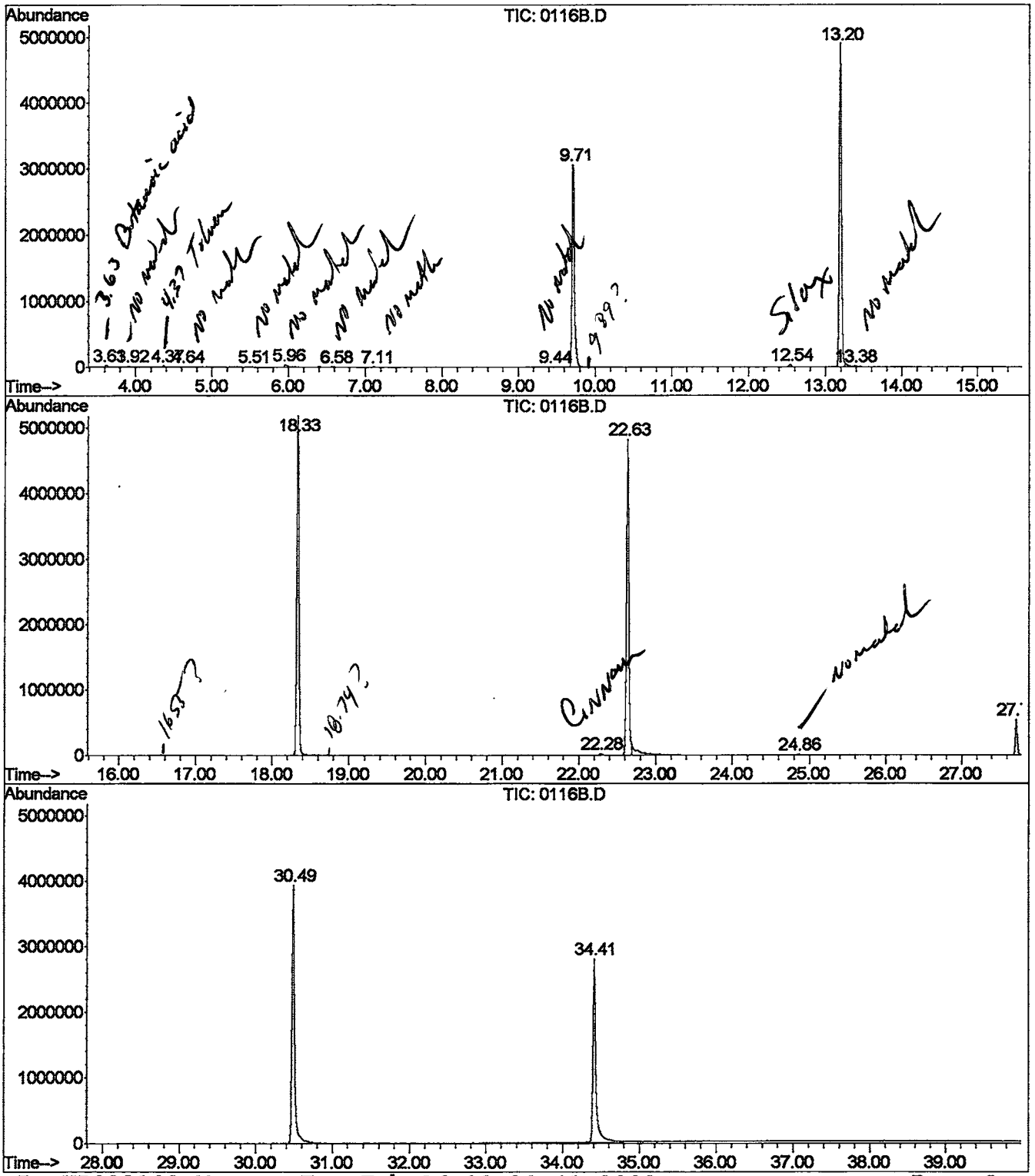
If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.611	19	21	27	rBV	21069	47094	0.46%	0.086%
2	3.916	43	48	51	rBV	13069	22985	0.23%	0.042%
3	4.367	85	88	95	rVB	17589	32682	0.32%	0.060%
4	4.638	111	112	117	rVB3	5831	13213	0.13%	0.024%
5	5.507	179	189	195	rBV2	5777	16697	0.16%	0.030%
6	5.959	225	229	243	rBV	29608	104049	1.02%	0.190%
7	6.580	279	284	287	rVV	5408	12132	0.12%	0.022%
8	7.110	325	331	335	rBV3	5761	15642	0.15%	0.029%
9	9.436	533	537	543	rVB2	3527	11734	0.12%	0.021%
10	9.707	557	561	581	rBV	3065175	6589498	64.70%	12.029%
11	12.540	807	812	817	rVB	34225	60901	0.60%	0.111%
12	13.195	865	870	875	rBV	4915499	9443091	92.72%	17.239%
13	13.376	885	886	891	rVB	9547	14273	0.14%	0.026%
14	18.332	1319	1325	1331	rBV	5184888	9978378	97.98%	18.216%
15	22.283	1671	1675	1681	rBV	21780	44563	0.44%	0.081%
16	22.633	1701	1706	1711	rBV2	4822382	10184228	100.00%	18.592%
17	24.857	1899	1903	1915	rVB	7201	21166	0.21%	0.039%
18	27.724	2153	2157	2163	rBV	542497	929199	9.12%	1.696%
19	30.490	2397	2402	2437	rBV	3929356	9576339	94.03%	17.482%
20	34.407	2743	2749	2783	rBV	2791071	7660702	75.22%	13.985%

Sum of corrected areas: 54778566

File : C:\MSDCHEM\1\5973N\02FEB25\0116B.D
Operator :
Acquired : 25 Feb 2002 11:14 am using AcqMethod HP020702
Instrument : Instrumen
Sample Name: lab blank 1/16
Misc Info : February 25, 2002
Vial Number: 1
Quant File :HP022102.RES (RTE Integrator)



Data File : C:\MSDCHEM\1\5973N\02FEB25\0116B.D Vial: 1
Acq On : 25 Feb 2002 11:14 am Operator:
Sample : lab blank 1/16 Inst : Instrumen
Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

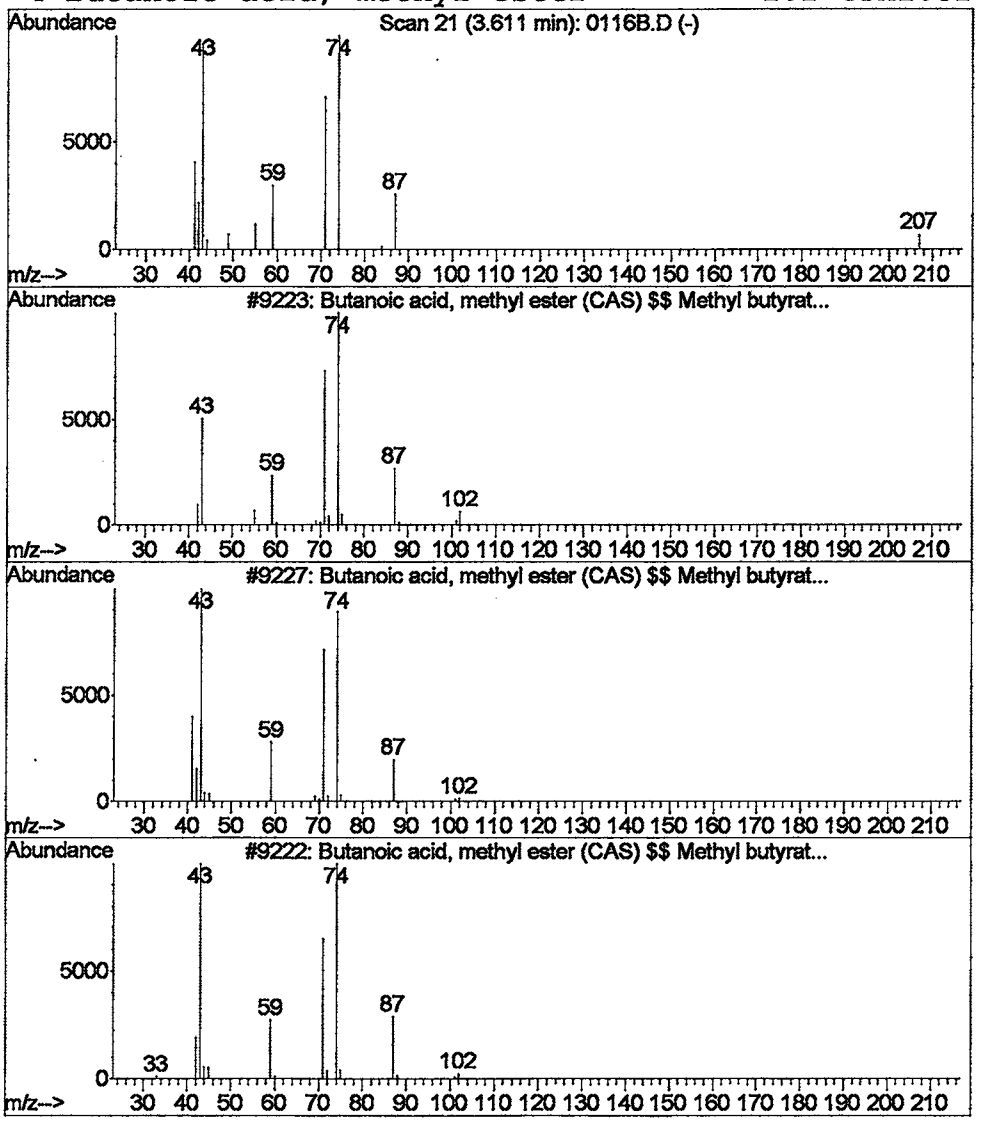
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

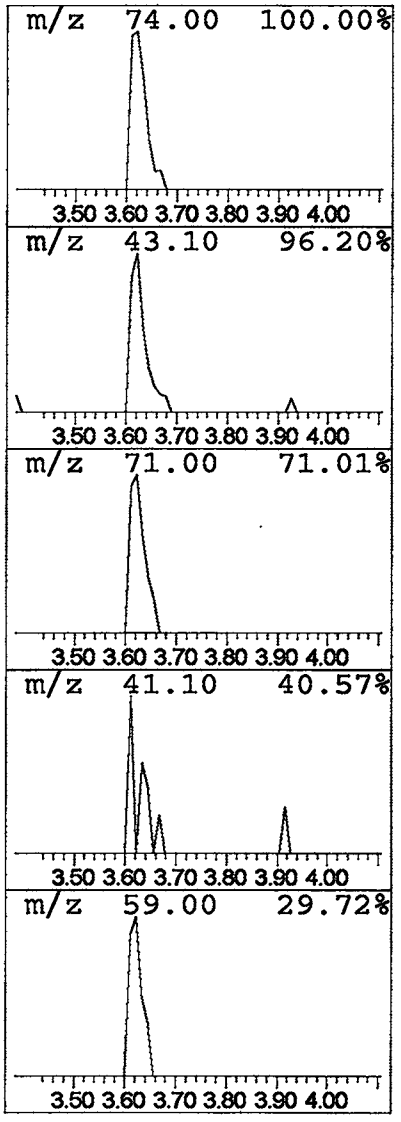
****

Peak Number 1 Butanoic acid, methyl ester... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.14 ug/L	47094	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
2	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
3	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
4	Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	83





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Acq On : 25 Feb 2002 11:14 am Operator:
Sample : lab blank 1/16 Inst : Instrumen
Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

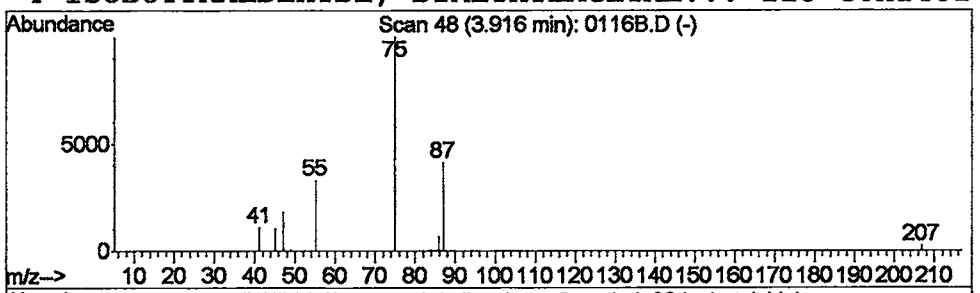
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

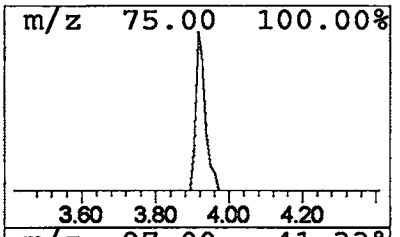
****

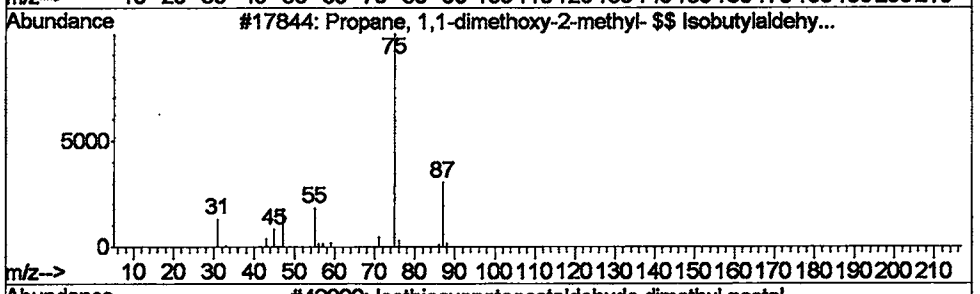
Peak Number 2 Propane, 1,1-dimethoxy-2-me... Concentration Rank 6

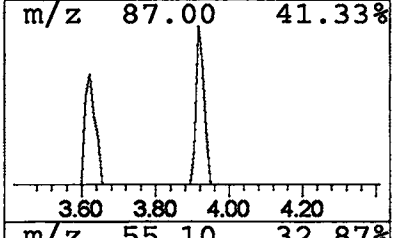
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.92	0.07 ug/L	22985	1,4-Dichlorobenzene-d4	9.72

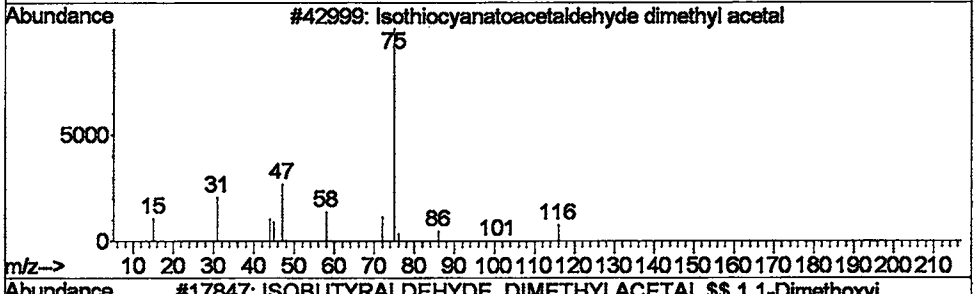
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,1-dimethoxy-2-methyl-...	118	C6H14O2	041632-89-7	9
2	Isothiocyanatoacetaldehyde dimet...	147	C5H9NO2S	075052-04-9	9
3	ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	9
4	ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	9

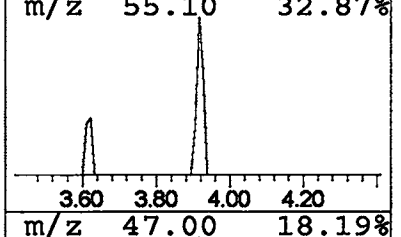


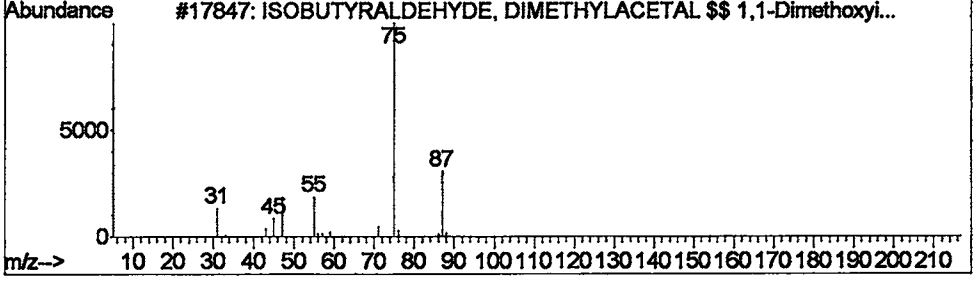


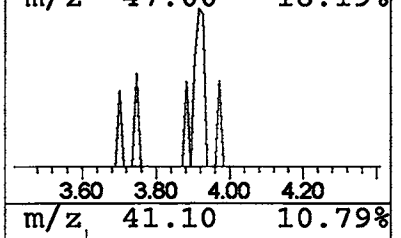


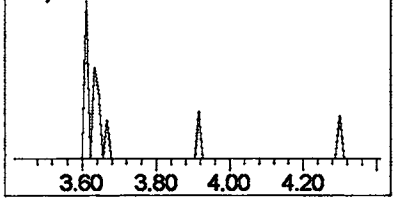












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Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

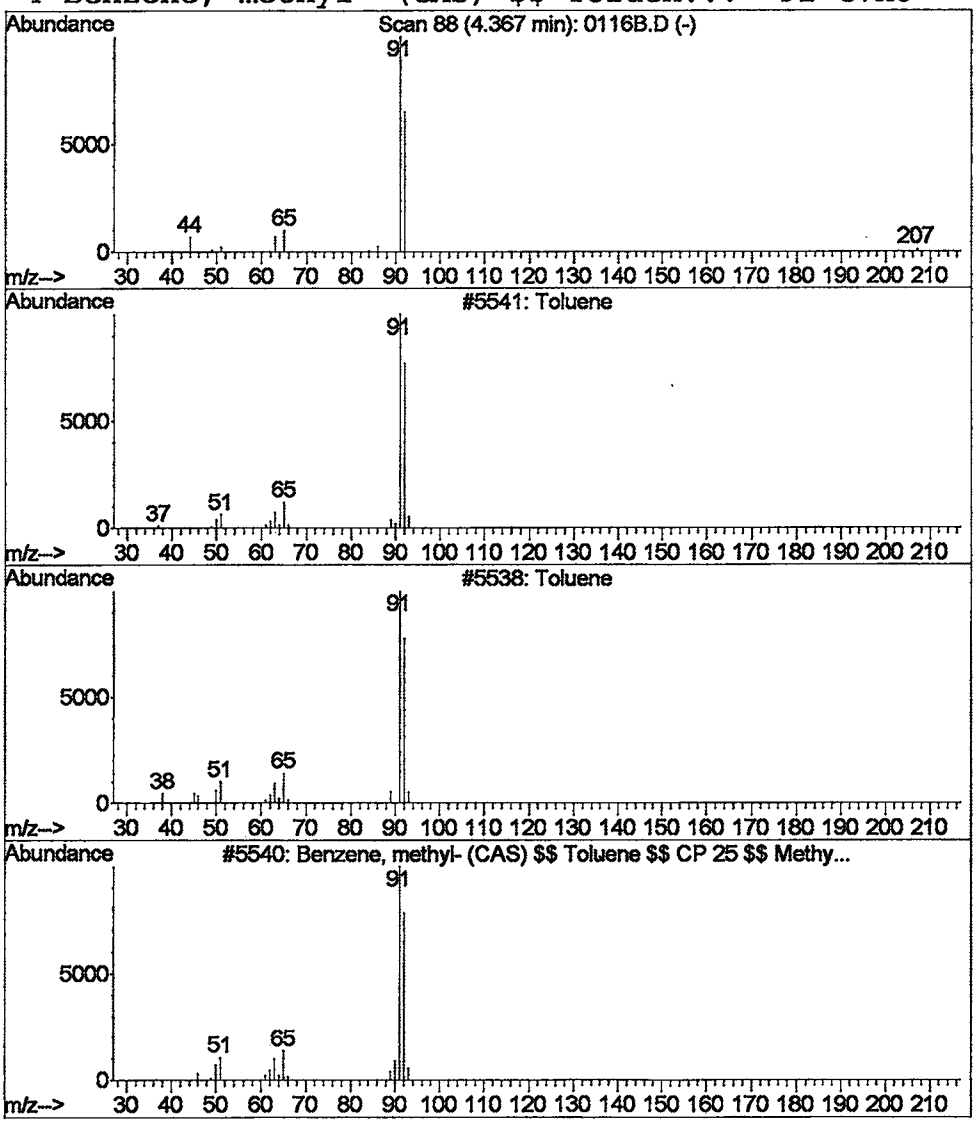
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

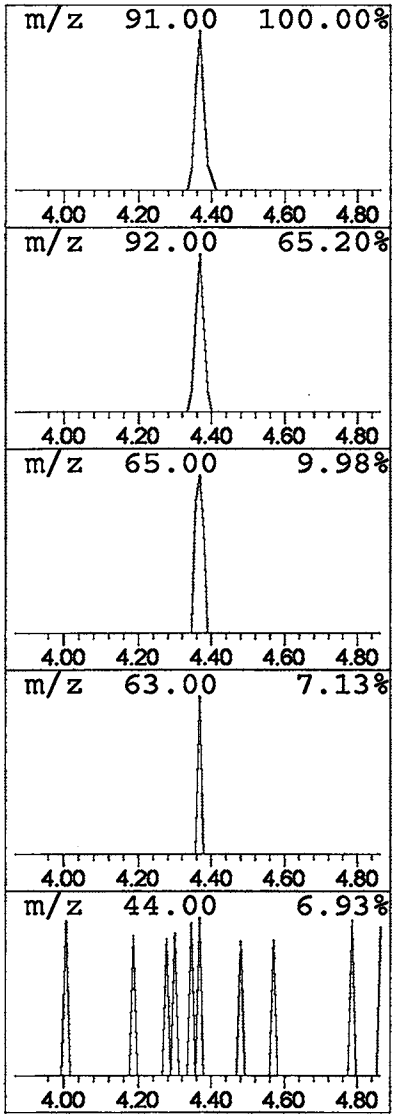
****

Peak Number 3 Toluene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	0.10 ug/L	32682	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	78
2	Toluene	92	C7H8	000108-88-3	78
3	Benzene, methyl- (CAS) $$ Toluen...	92	C7H8	000108-88-3	78
4	Benzene, methyl- (CAS) $$ Toluen...	92	C7H8	000108-88-3	64





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Acq On : 25 Feb 2002 11:14 am Operator:
Sample : lab blank 1/16 Inst : Instrumen
Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

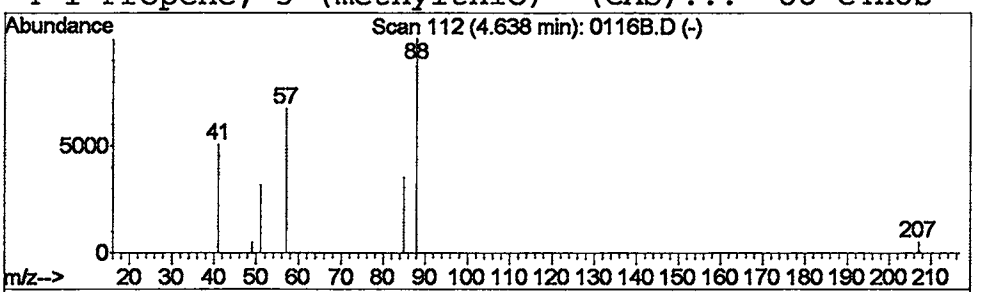
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

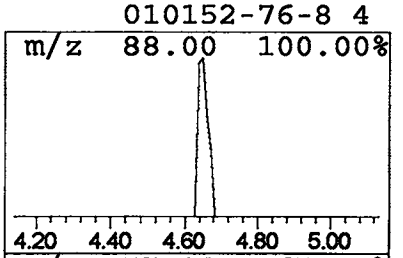
****

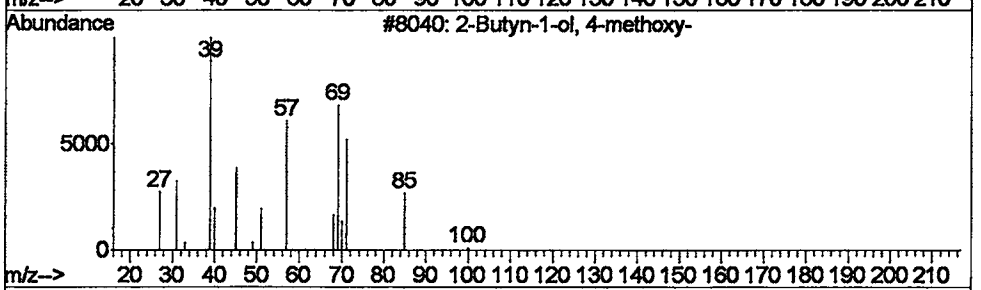
Peak Number 4 2-Butyn-1-ol, 4-methoxy- Concentration Rank 10

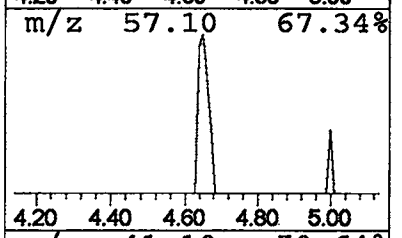
R.T.	EstConc	Area	Relative to ISTD	R.T.
4.64	0.04 ug/L	13213	1,4-Dichlorobenzene-d4	9.72

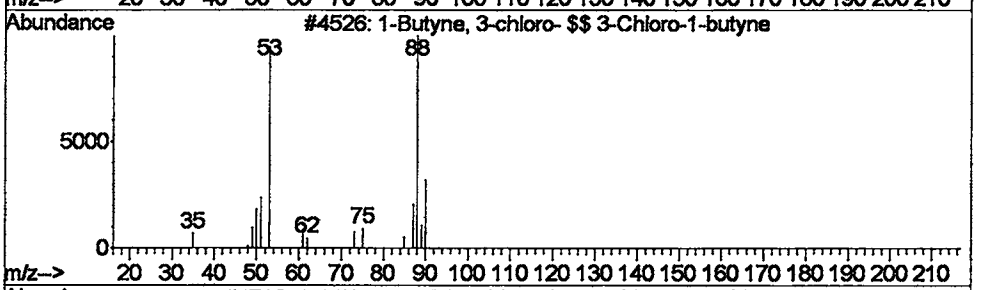
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butyn-1-ol, 4-methoxy-	100	C5H8O2	018857-03-9	9
2	1-Butyne, 3-chloro- $$ 3-Chloro-...	88	C4H5Cl	021020-24-6	4
3	1,4-Dioxane (CAS) $$ p-Dioxane $...	88	C4H8O2	000123-91-1	4
4	1-Propene, 3-(methylthio)- (CAS)...	88	C4H8S	010152-76-8	4

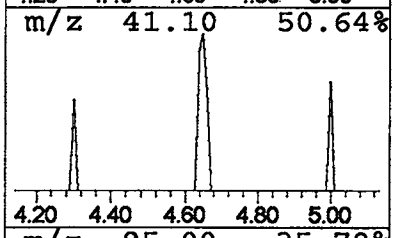


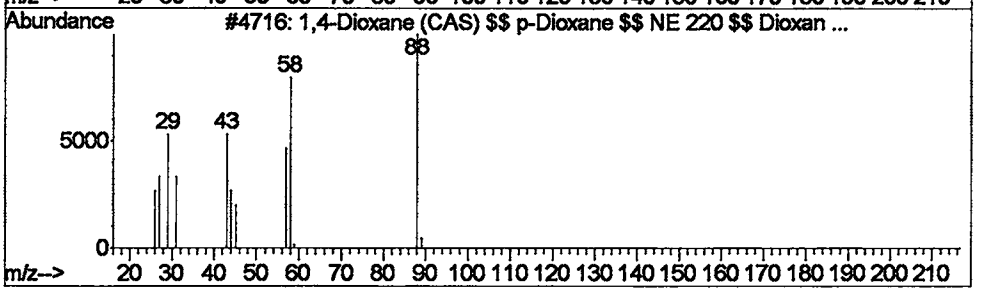


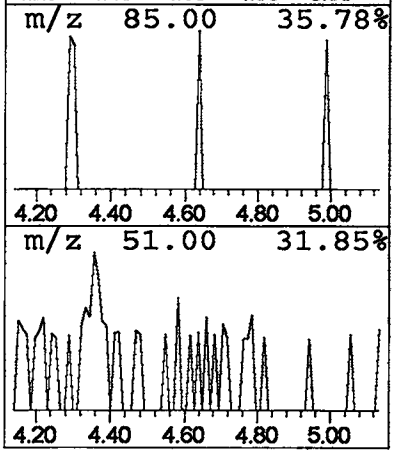












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Acq On : 25 Feb 2002 11:14 am Operator:
Sample : lab blank 1/16 Inst : Instrumen
Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

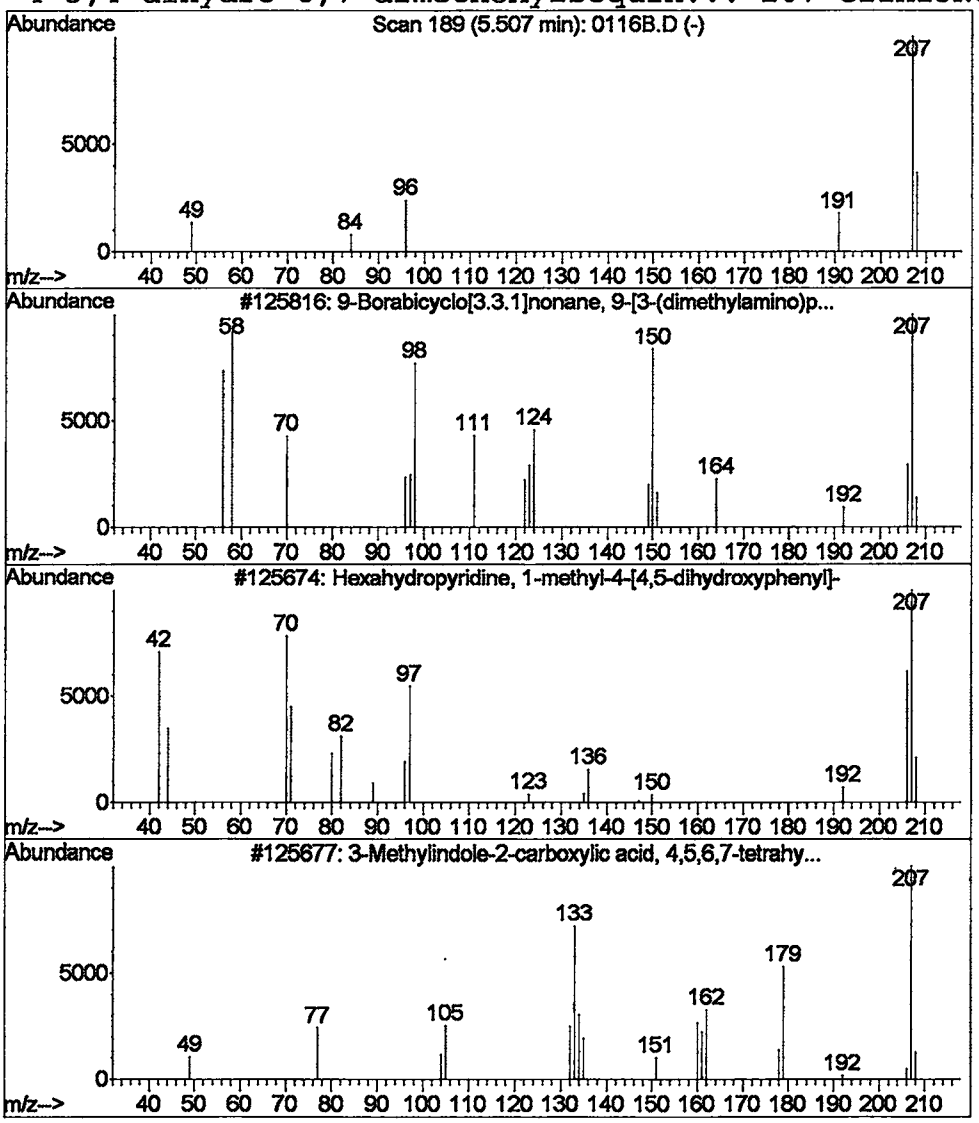
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

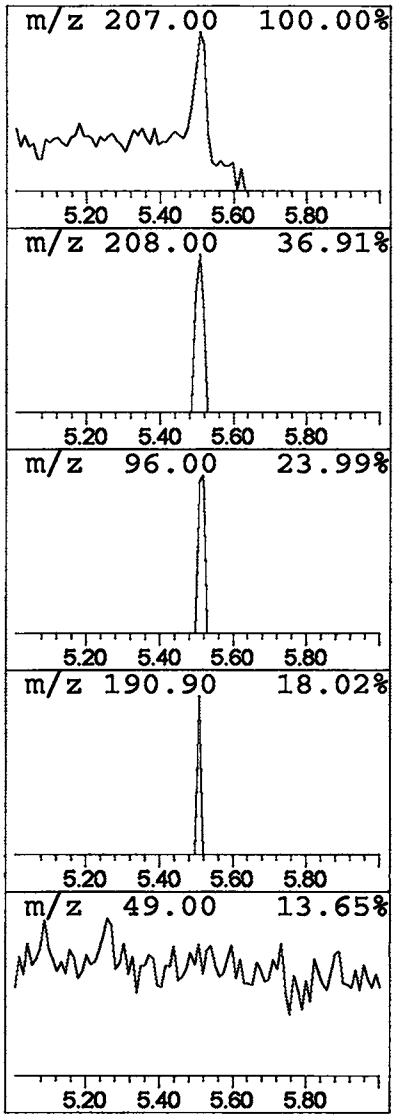
****

Peak Number 5 9-Borabicyclo[3.3.1]nonane,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.51	0.05 ug/L	16697	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Borabicyclo[3.3.1]nonane, 9-[3...	207	C13H26BN	000000-00-0	5
2	Hexahydropyridine, 1-methyl-4-[4...	207	C12H17NO2	094427-47-1	5
3	3-Methylindole-2-carboxylic acid...	207	C12H17NO2	000000-00-0	5
4	3,4-dihydro-6,7-dimethoxyisoquin...	207	C11H13NO3	084122-10-1	5





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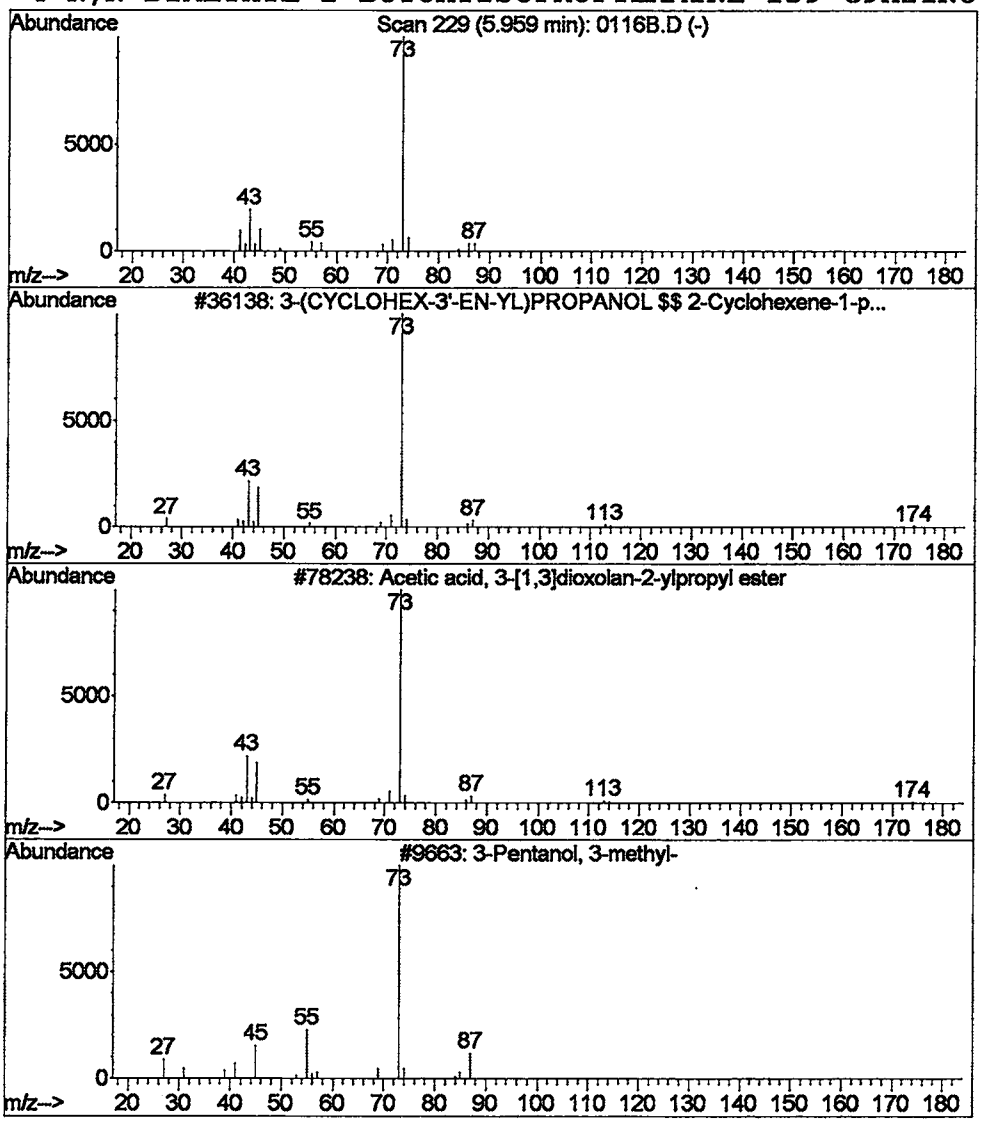
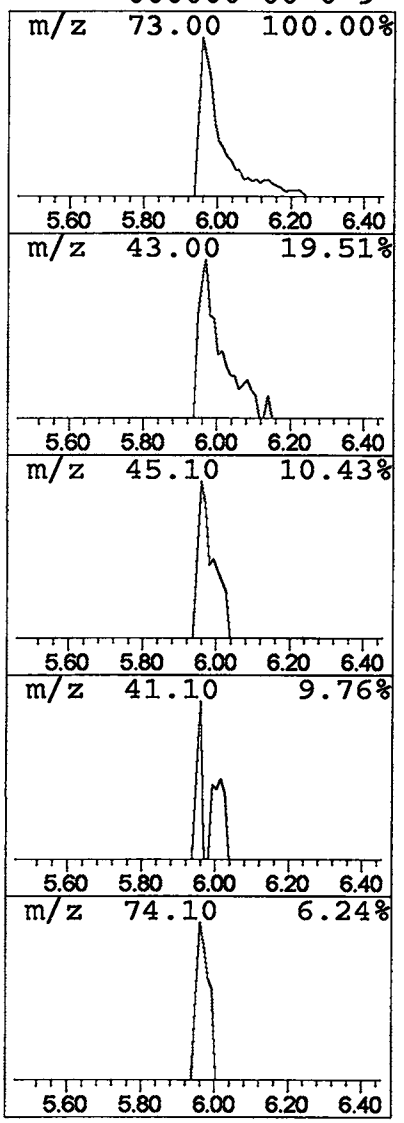
****

Peak Number 6 3-(CYCLOHEX-3'-EN-YL)PROPAN... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.96	0.32 ug/L	104049	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-(CYCLOHEX-3'-EN-YL)PROPANOL $$...	174	C8H14O4	015745-87-6	33
2	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	33
3	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	33
4	N,N-DIMETHYL-2-BUTOXYISOPROPYLAMINE	159	C9H21NO	000000-00-0	9





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Acq On    : 25 Feb 2002  11:14 am                    Operator:
Sample    : lab blank 1/16                           Inst    : Instrumen
Misc      : February 25, 2002                        Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title        : 508 scan method
Library      : C:\DATABASE\WILEY7N.L
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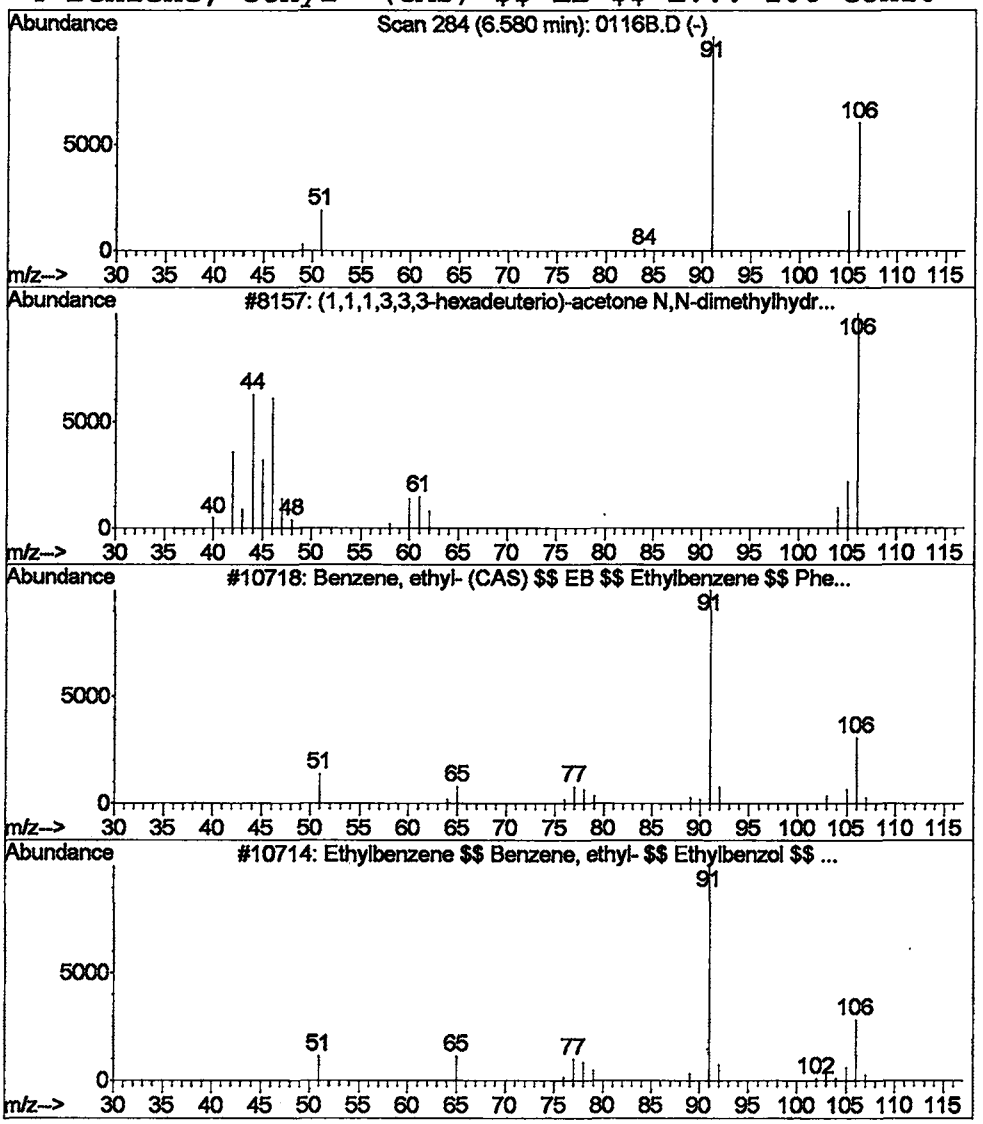
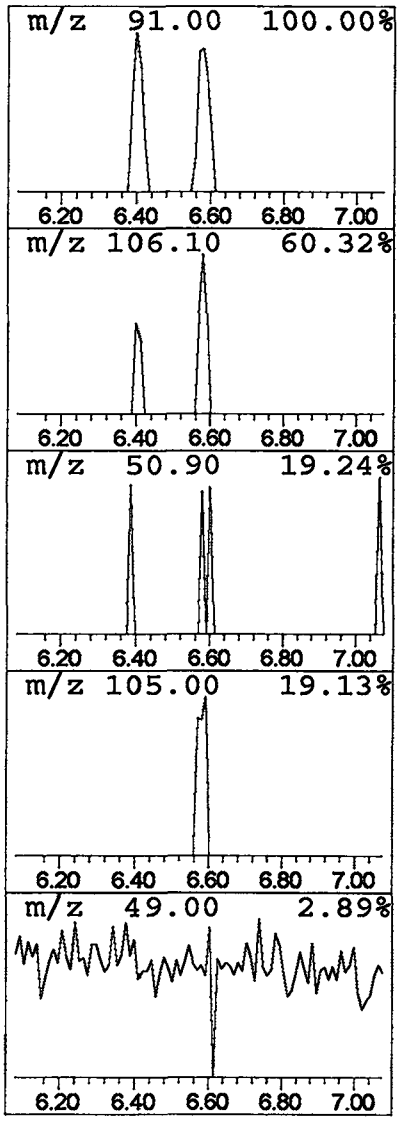
Peak Number 7 (1,1,1,3,3,3-hexadeuterio)-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	0.04 ug/L	12132	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1,1,1,3,3,3-hexadeuterio)-aceto...	106	C5H6D6N2	123499-00-3	4
2	Benzene, ethyl- (CAS) $$ EB $$ E...	106	C8H10	000100-41-4	4
3	Ethylbenzene $$ Benzene, ethyl- ...	106	C8H10	000100-41-4	4
4	Benzene, ethyl- (CAS) $$ EB $$ E...	106	C8H10	000100-41-4	4





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MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

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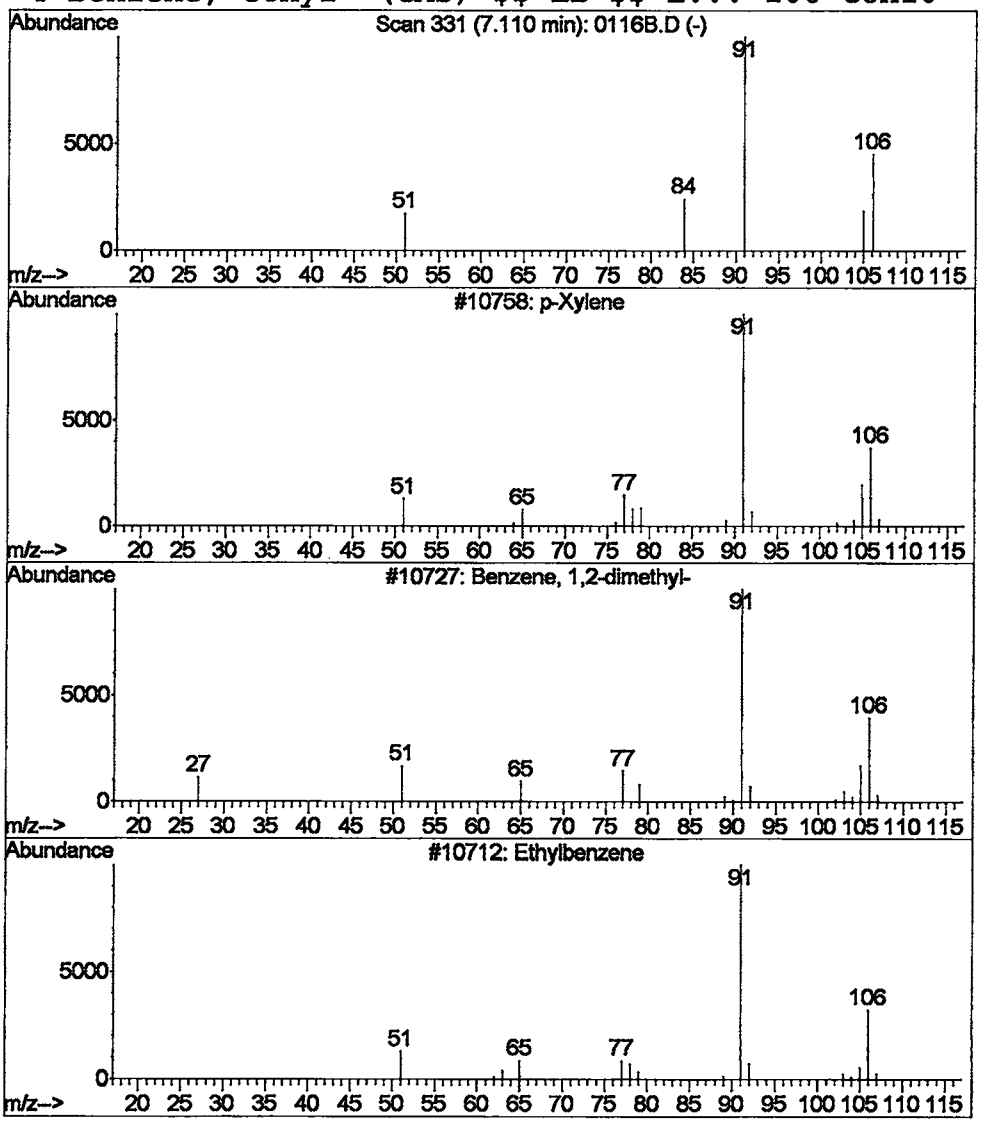
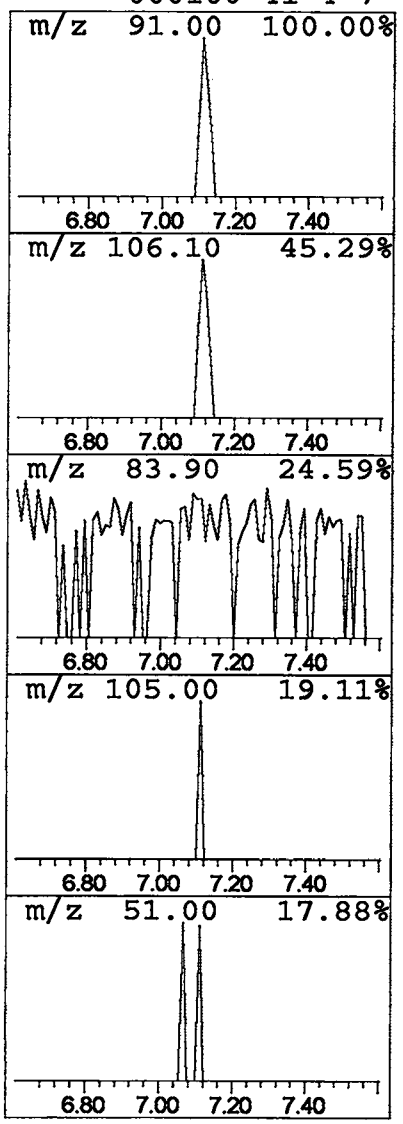
Peak Number 8 p-Xylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	0.05 ug/L	15642	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Xylene	106	C8H10	000106-42-3	45
2	Benzene, 1,2-dimethyl-	106	C8H10	000095-47-6	45
3	Ethylbenzene	106	C8H10	000100-41-4	7
4	Benzene, ethyl- (CAS) $$ EB $$ E...	106	C8H10	000100-41-4	7

NO GOOD MATCH





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MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

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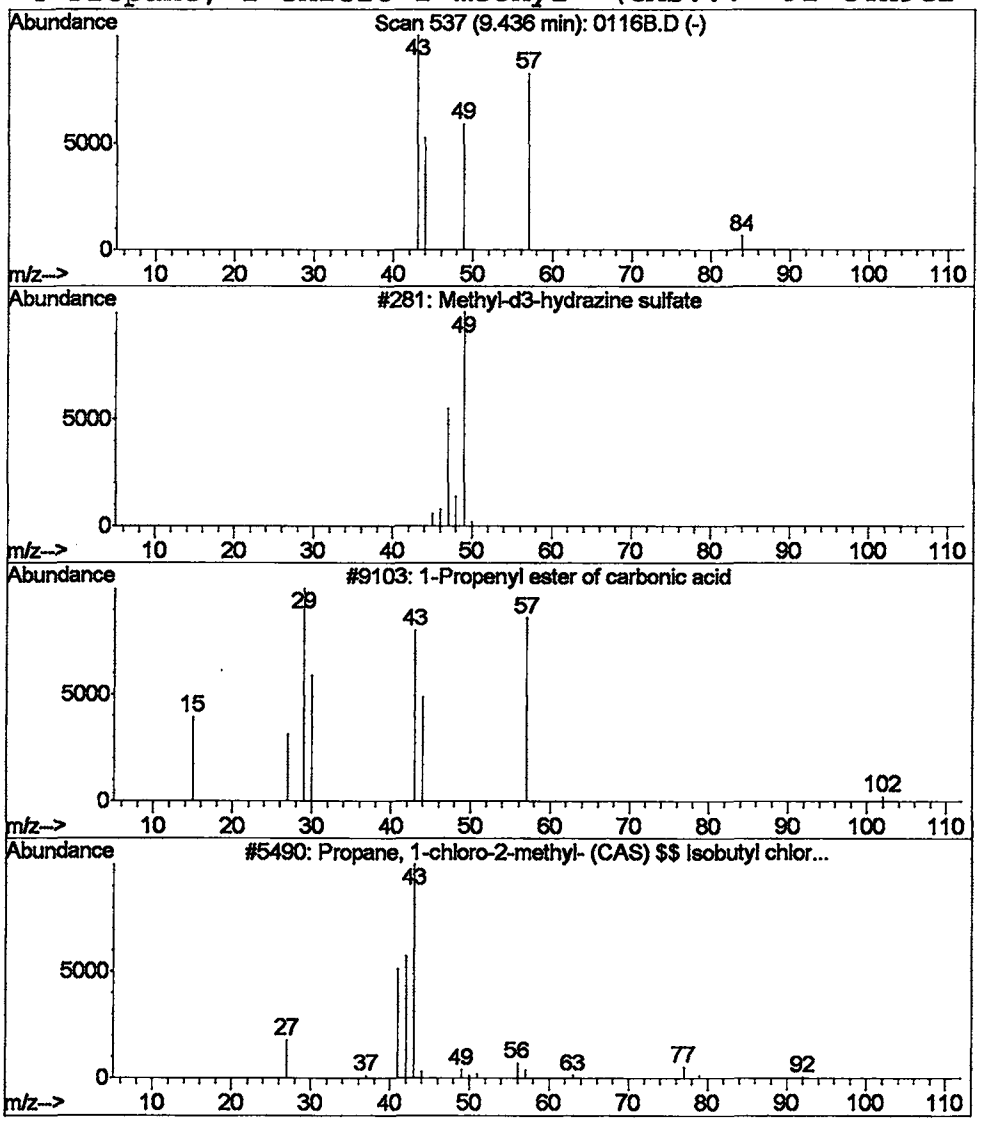
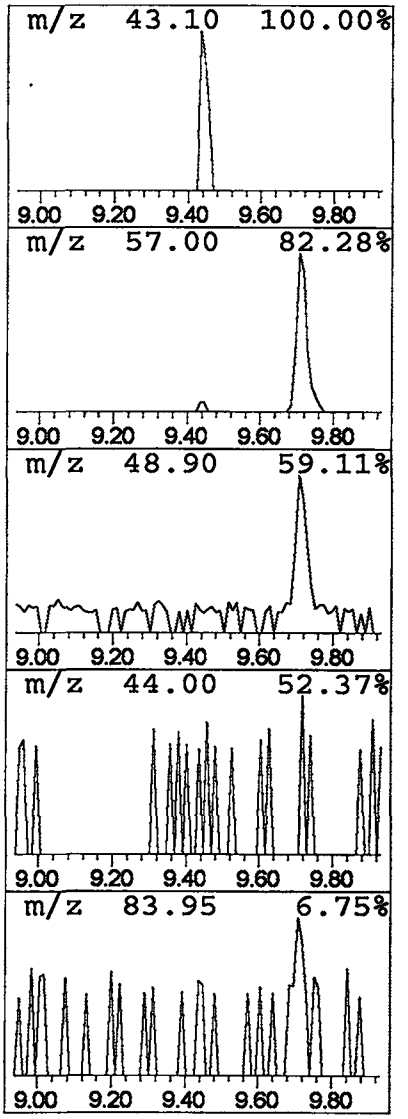
Peak Number 9 Methyl-d3-hydrazine sulfate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	0.04 ug/L	11734	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl-d3-hydrazine sulfate	49	CH3D3N2	000000-00-0	3
2	1-Propenyl ester of carbonic acid	102	C4H6O3	000000-00-0	2
3	Propane, 1-chloro-2-methyl- (CAS...	92	C4H9Cl	000513-36-0	2
4	Propane, 1-chloro-2-methyl- (CAS...	92	C4H9Cl	000513-36-0	2





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Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

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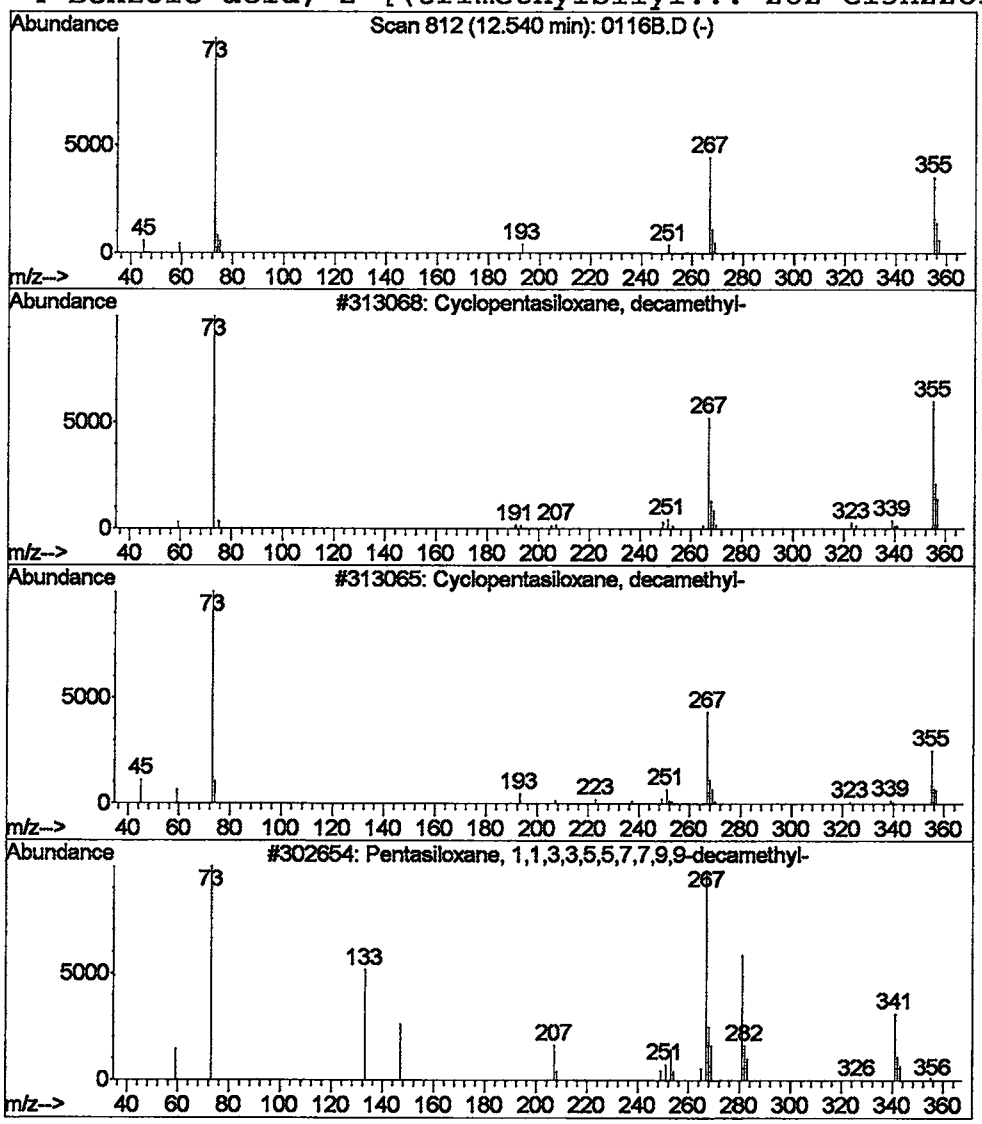
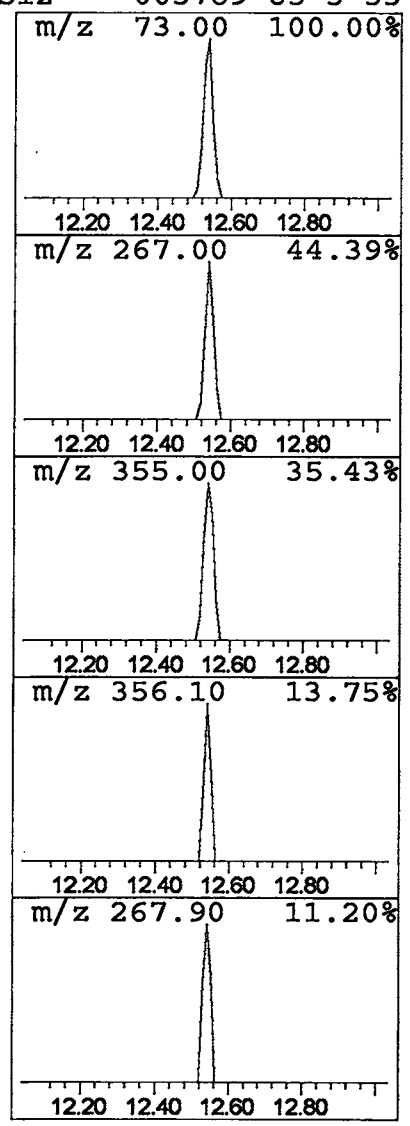
Peak Number 10 Cyclopentasiloxane, decamet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.13 ug/L	60901	Naphthalene-d8	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	50
2	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	38
3	Pentasiloxane, 1,1,3,3,5,5,7,7,9...	356	C10H32O4Si5	000995-83-5	36
4	Benzoic acid, 2-[(trimethylsilyl...	282	C13H22O3Si2	003789-85-3	33





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MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

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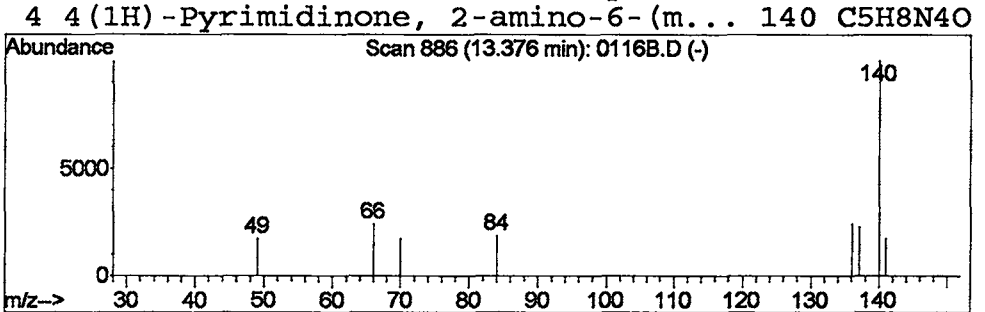
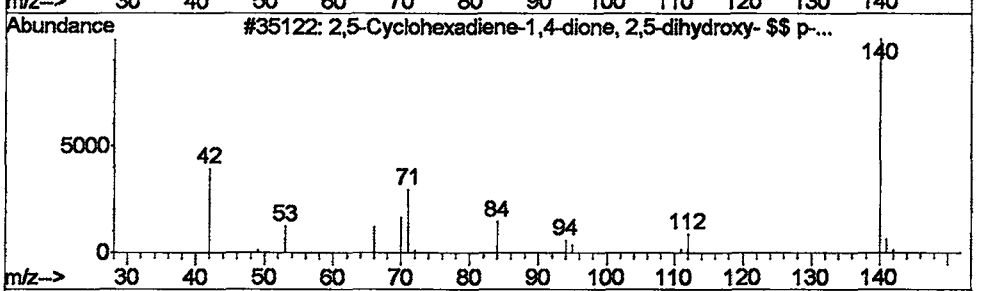
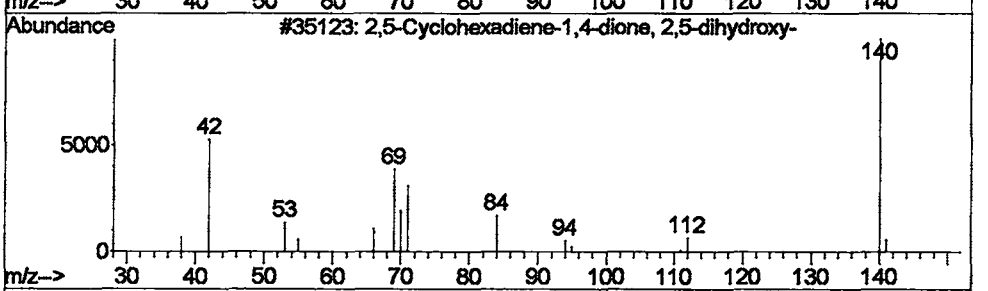
Peak Number 11 2,5-Cyclohexadiene-1,4-dion... Concentration Rank 13

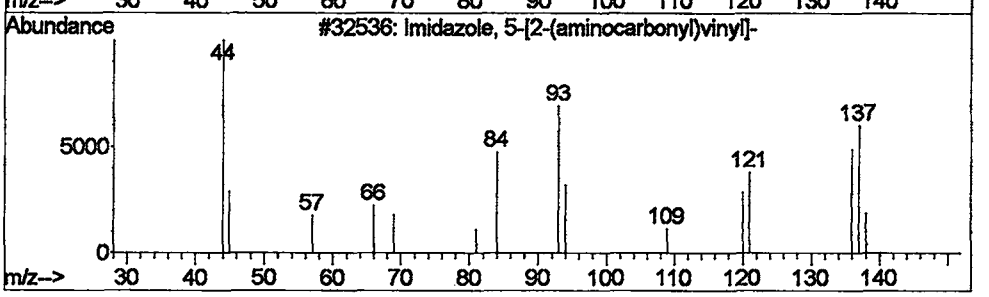
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	0.03 ug/L	14273	Naphthalene-d8	13.20

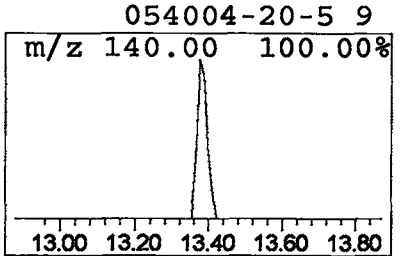
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	9
2	2,5-Cyclohexadiene-1,4-dione, 2,...	140	C6H4O4	000615-94-1	9
3	Imidazole, 5-[2-(aminocarbonyl)v...	137	C6H7N3O	135200-64-5	9
4	4(1H)-Pyrimidinone, 2-amino-6-(m...	140	C5H8N4O	054004-20-5	9

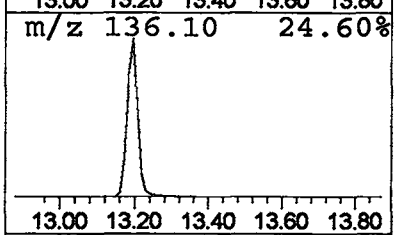


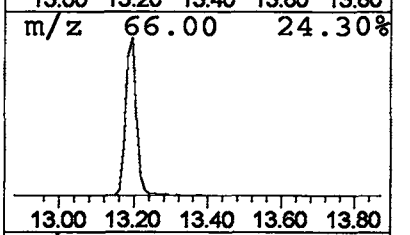


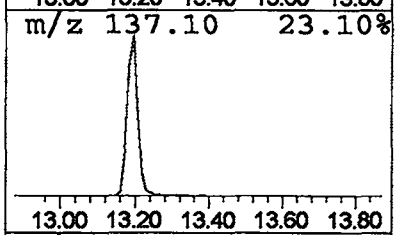


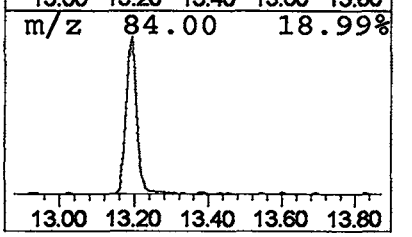












Data File : C:\MSDCHEM\1\5973N\02FEB25\0116B.D Vial: 1
Acq On : 25 Feb 2002 11:14 am Operator:
Sample : lab blank 1/16 Inst : Instrumen
Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

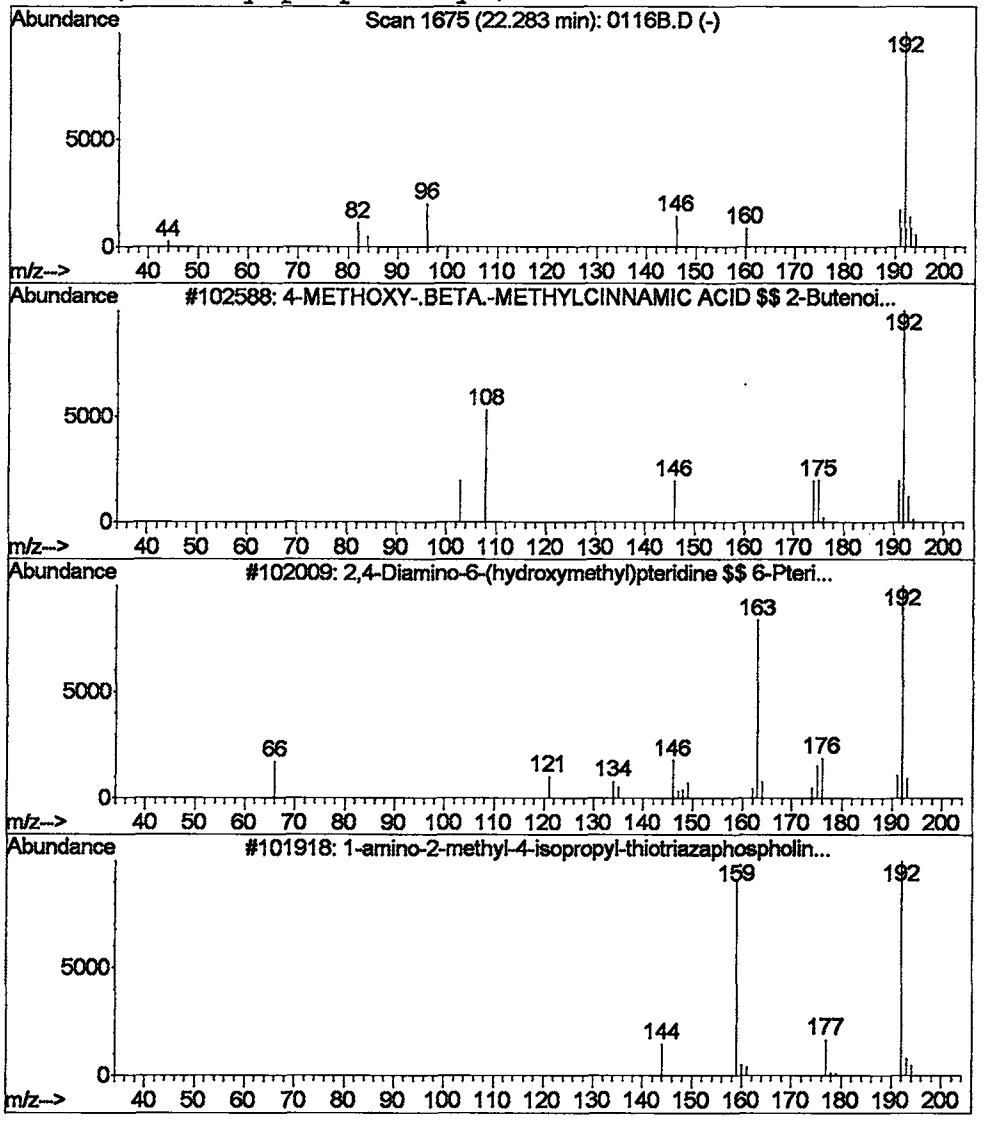
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

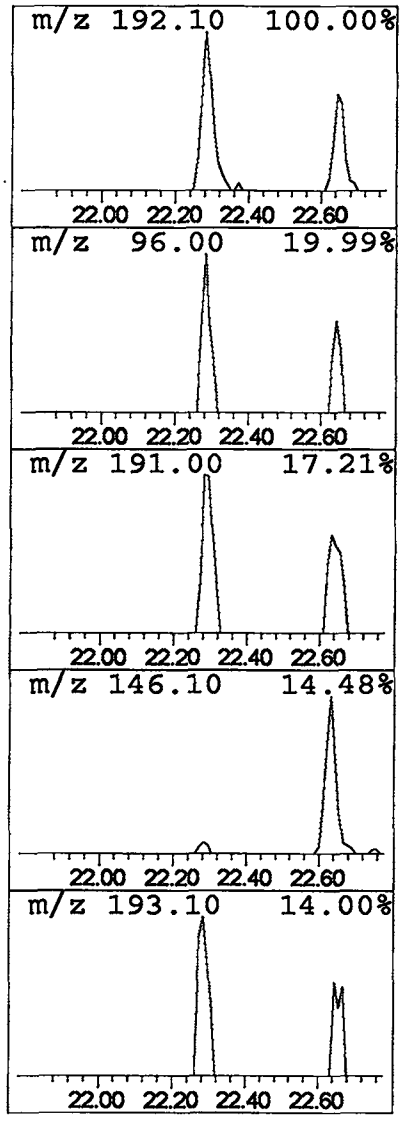
****

Peak Number 12 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.09 ug/L	44563	Phenanthrene-d10	22.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-METHOXY-.BETA.-METHYLCINNAMIC ...	192	C11H12O3	019169-94-9	83
2	2,4-Diamino-6-(hydroxymethyl)pte...	192	C7H8N6O	000945-24-4	50
3	1-amino-2-methyl-4-isopropyl-thi...	192	C5H13N4PS	107264-54-0	50
4	N-(2-methylprop-2-enyl)-2-nitrob...	192	C10H12N2O2	000000-00-0	50





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Misc : February 25, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

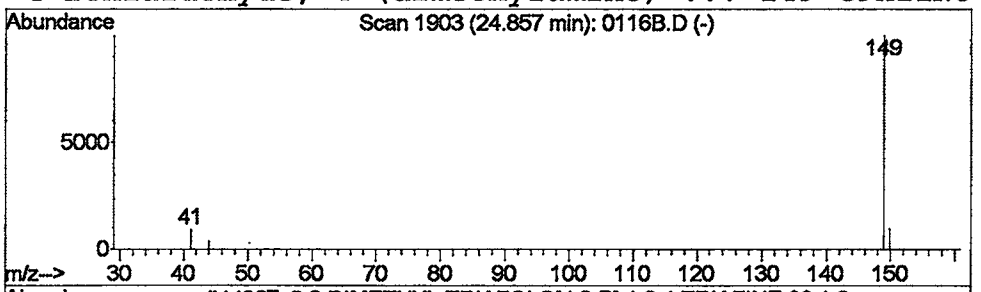
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

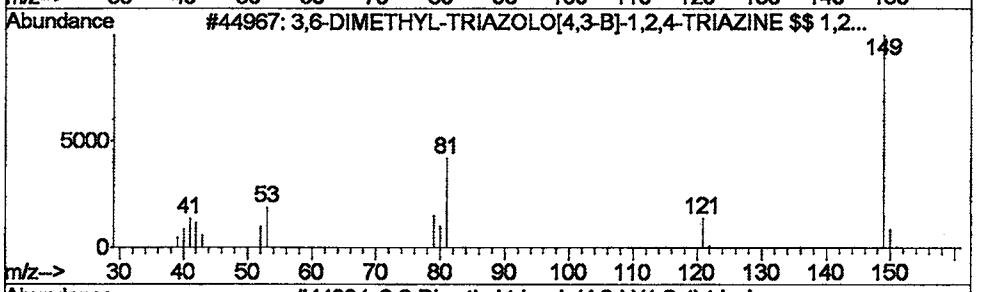
****

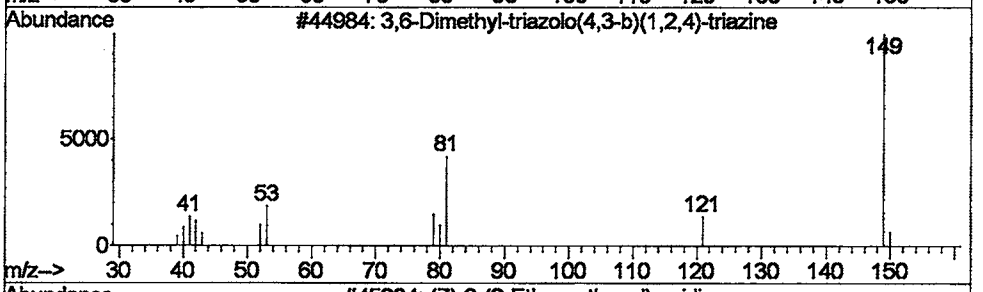
Peak Number 13 3,6-DIMETHYL-TRIAZOLO[4,3-B... Concentration Rank 9

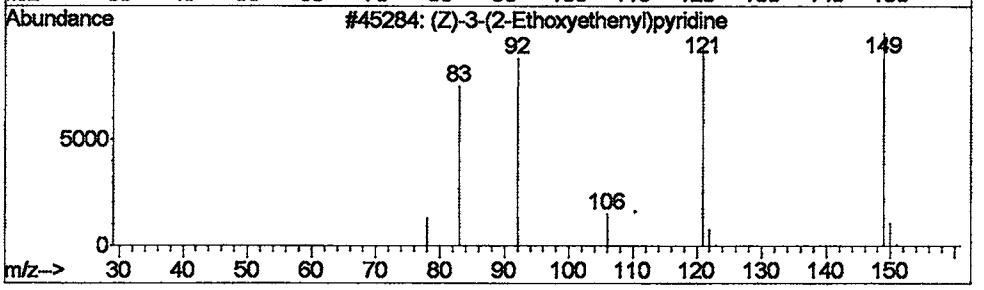
R.T.	EstConc	Area	Relative to ISTD	R.T.
24.86	0.04 ug/L	21166	Phenanthrene-d10	22.63

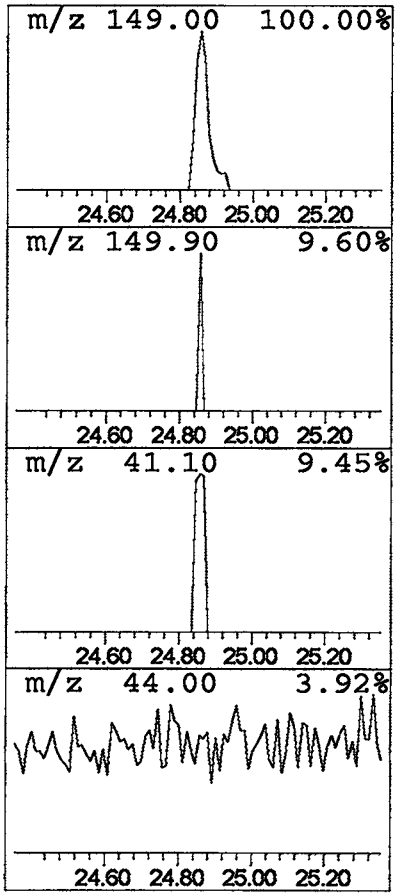
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6-DIMETHYL-TRIAZOLO[4,3-B]-1,2...	149	C6H7N5	061139-71-7	5
2	3,6-Dimethyl-triazolo(4,3-b)(1,2...	149	C6H7N5	000000-00-0	5
3	(Z)-3-(2-Ethoxyethenyl)pyridine	149	C9H11NO	000000-00-0	4
4	Benzaldehyde, 4-(dimethylamino)-...	149	C9H11NO	000100-10-7	4











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Name: lab blank 1/16
Misc: February 25, 2002
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Title: 508 scan method
Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--- #	RT	Resp	Conc
Butanoic acid, me...	3.61	0.1	ug/L	47094	1	9.72	6589500	20.0
Propane, 1,1-dime...	3.92	0.1	ug/L	22985	1	9.72	6589500	20.0
Toluene	4.37	0.1	ug/L	32682	1	9.72	6589500	20.0
2-Butyn-1-ol, 4-m...	4.64	0.0	ug/L	13213	1	9.72	6589500	20.0
9-Borabicyclo[3.3...	5.51	0.1	ug/L	16697	1	9.72	6589500	20.0
3-(CYCLOHEX-3'-EN...	5.96	0.3	ug/L	104049	1	9.72	6589500	20.0
(1,1,1,3,3,3-hexa...	6.58	0.0	ug/L	12132	1	9.72	6589500	20.0
p-Xylene	7.11	0.0	ug/L	15642	1	9.72	6589500	20.0
Methyl-d3-hydrazi...	9.44	0.0	ug/L	11734	1	9.72	6589500	20.0
Cyclopentasiloxan...	12.54	0.1	ug/L	60901	2	13.20	9443090	20.0
2,5-Cyclohexadien...	13.38	0.0	ug/L	14273	2	13.20	9443090	20.0
4-METHOXY-.BETA.-...	22.28	0.1	ug/L	44563	4	22.63	10184200	20.0
3,6-DIMETHYL-TRIA...	24.86	0.0	ug/L	21166	4	22.63	10184200	20.0