User: Baglioni, Walter
Westchester Dept. of Labs & Research
Date: 01/31/2002 Time: 14:17:25

Sample: AE00555
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 1/29/02 14:16 Ended: 1/31/02 14:16 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	Hexachlorobutadiene		< 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-Diphenylhydraz		< 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

NYC 9/11 Public Portal Document

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Jan 30 08:26:04 2002 Quant Results File: SCAN508.RES

Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Last Update : Fri Jan 11 13:20:07 2002
Response via : Initial Calibration
DataAcq Meth : SCAN508

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.71	152	1513741	20.00	ug/L	0.00
10) Naphthalene-d8	13.19	136	6490883	20.00	ug/L	-0.02
17) Acenaphthene-d10	18.34	164	3352807	20.00	ug/L	-0.01
29) Phenanthrene-d10	22.65	188	5496787	20.00	ug/L	-0.02
38) Chrysene-d12	30.50	240	4687760	20.00	ug/L	-0.01
44) Perylene-d12	34.42	264	3369095	20.00	ug/L	0.00
System Monitoring Compounds						
37) 2,4-DDD	27.73	235	322914	3.72	ug/L	0.00
Spiked Amount 5.000			Recovery = 74.40%			
Target Compounds						Qvalue
20) Dimethylphthalate	18.34	163	534588	2.41	ug/L #	56
21) 2,6-Dinitrotoluene	18.34	165	422260	9.78	ug/L #	17

(#) = qualifier out of range (m) = manual integration
0109B.D SCAN508.M Wed Jan 30 08:26:05 2002 Page 1

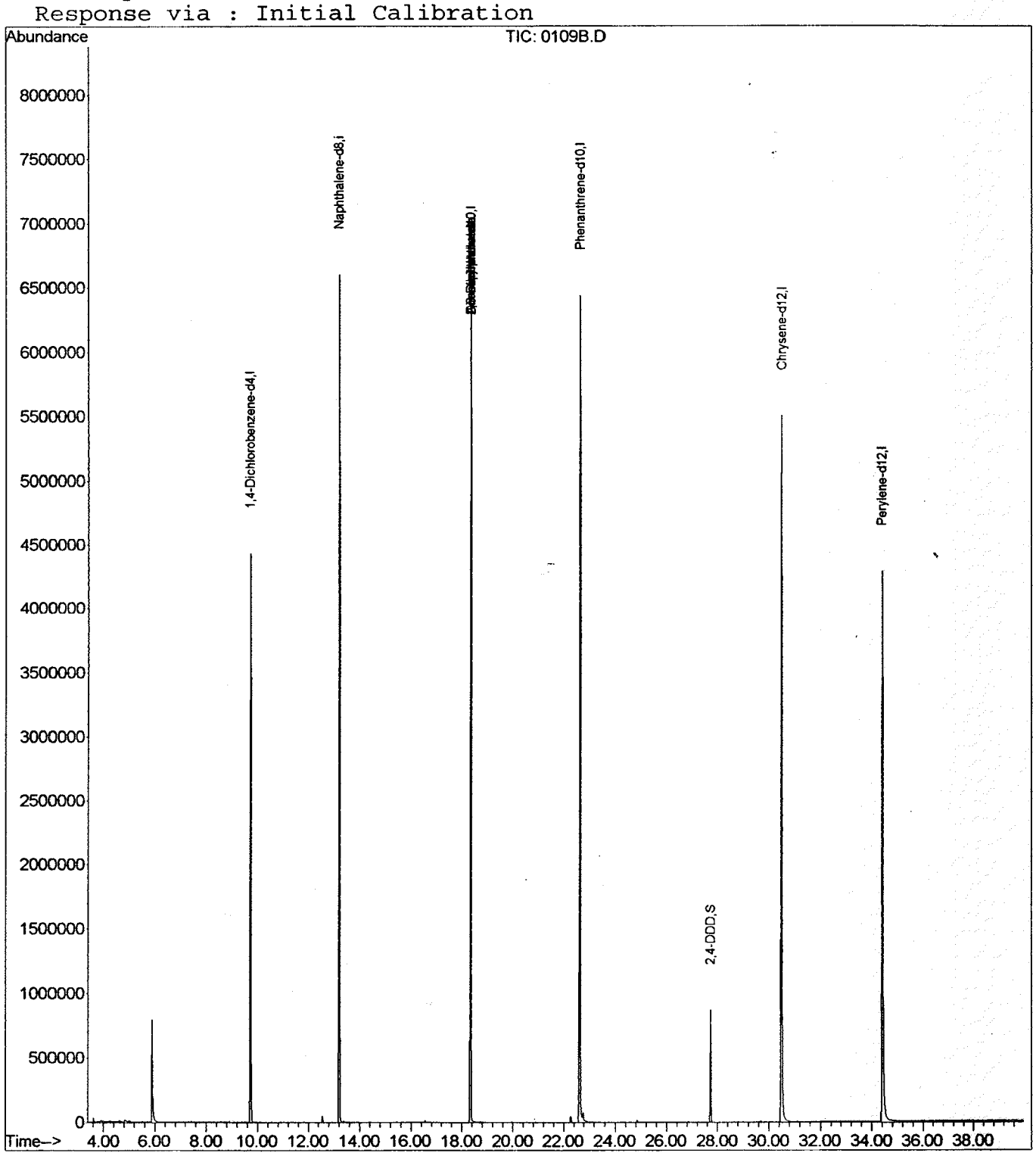
NYC-WTC_000051513

NYC 9/11 Public Portal Document

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Jan 30 8:26 2002 Quant Results File: SCAN508.RE

Method : C:\MSDCHEM\1\METHODS\METHODS\SCAN508.M (RTE Integrator)
Title : 508 scan method
Last Update : Fri Jan 11 13:20:07 2002
Response via : Initial Calibration



NYC-WTC_000051514

NYC 9/11 Public Portal Document

LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\MSDCHEM\1\METHODS\METHODS\SCAN508.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	3.602	17	20	25	rBV	33532	51376	0.35%	0.066%
2	3.910	45	47	51	rVB	12416	21080	0.14%	0.027%
3	4.253	73	77	85	rVB4	9470	28388	0.19%	0.036%
4	4.630	107	110	119	rVB2	15943	36317	0.25%	0.046%
5	4.847	125	129	139	rBV4	19530	76299	0.52%	0.097%
6	5.006	139	143	149	rVB3	11546	27816	0.19%	0.035%
7	5.885	215	220	243	rVB	796312	1645943	11.25%	2.100%
8	9.710	551	555	561	rBV	4430639	8739383	59.75%	11.151%
9	12.542	799	803	807	rVB2	48663	81429	0.56%	0.104%
10	13.192	855	860	865	rBV	6602695	12499879	85.47%	15.949%
11	16.549	1149	1154	1159	rVB	12458	22910	0.16%	0.029%
12	18.341	1305	1311	1315	rBV	6985472	13567485	92.77%	17.312%
13	22.280	1653	1656	1661	rBV2	47019	89849	0.61%	0.115%
14	22.645	1681	1688	1693	rBV	6440076	14625494	100.00%	18.662%
15	22.771	1697	1699	1717	rVB	75416	209940	1.44%	0.268%
16	24.849	1875	1881	1885	rVV	17308	35251	0.24%	0.045%
17	27.726	2127	2133	2139	rVB	865786	1566927	10.71%	1.999%
18	29.678	2299	2304	2311	rVV	7082	24567	0.17%	0.031%
19	30.500	2369	2376	2381	rBV	5500927	13557484	92.70%	17.299%
20	34.416	2713	2719	2747	rBV	4286358	11463824	78.38%	14.628%

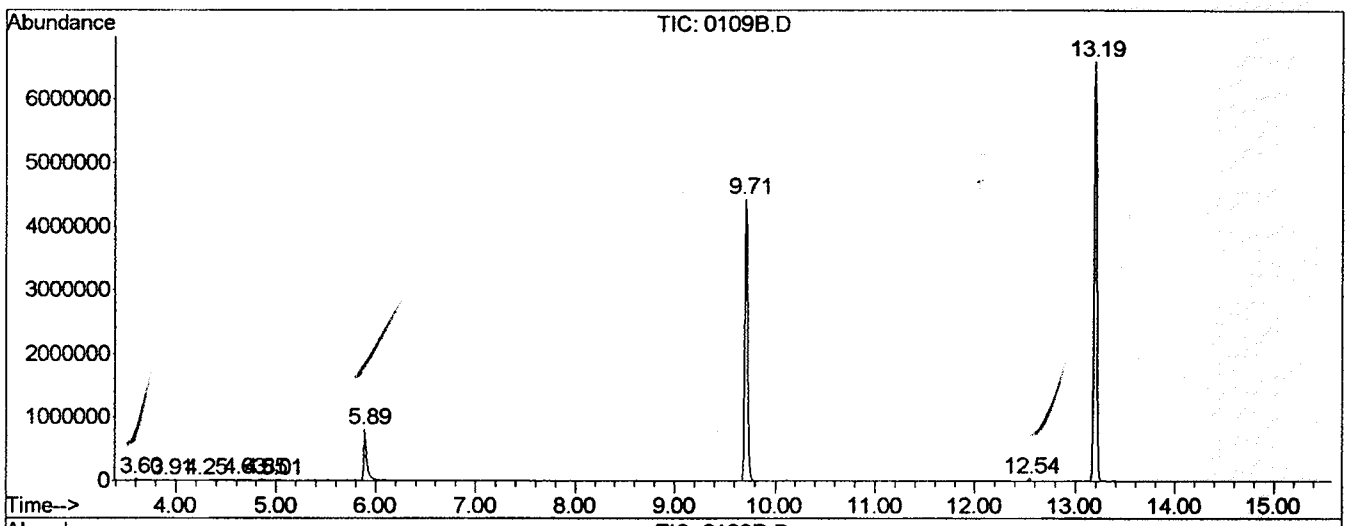
Sum of corrected areas: 78371641

NYC-WTC_000051515

NYC 9/11 Public Portal Document

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D
Operator :
Acquired : 29 Jan 2002 10:31 am using AcqMethod SCAN508
Instrument : Instrumen
Sample Name: lab blank 1/9
Misc Info : January 29, 2002
Vial Number: 1
Quant File :SCAN508.RES (RTE Integrator)



0109B.D SCAN508.M Wed Jan 30 09:02:32 2002 Page 2

NYC-WTC_000051516

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

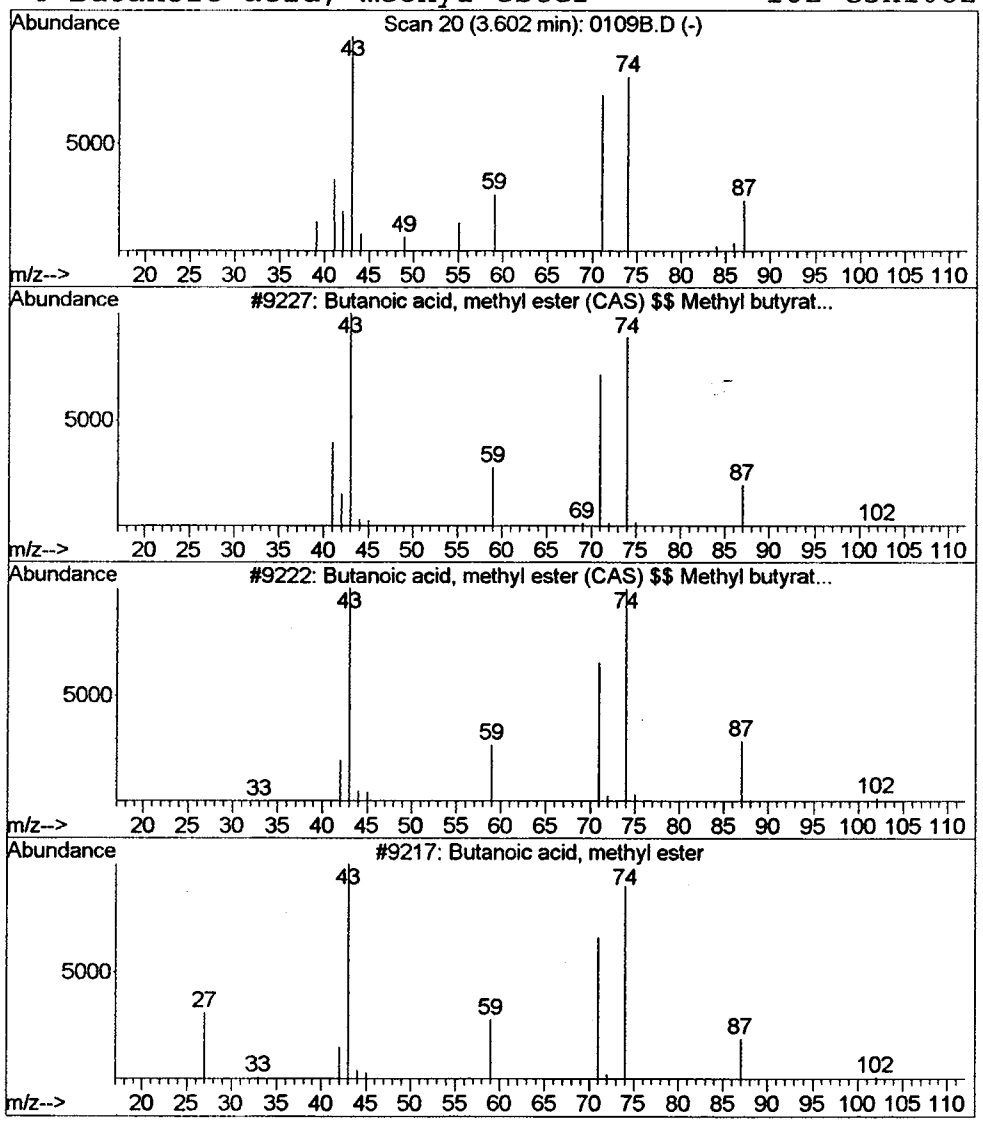
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

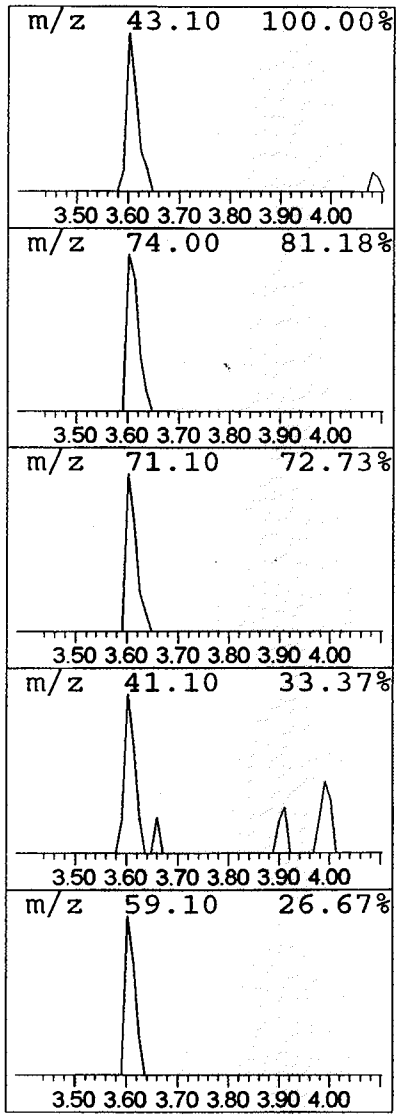
****

Peak Number 1 Butanoic acid, methyl ester... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.60	0.12 ug/L	51376	1,4-Dichlorobenzene-d4	9.71

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	102	C5H10O2	000623-42-7	83
4	Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	78





NYC-WTC_000051517

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

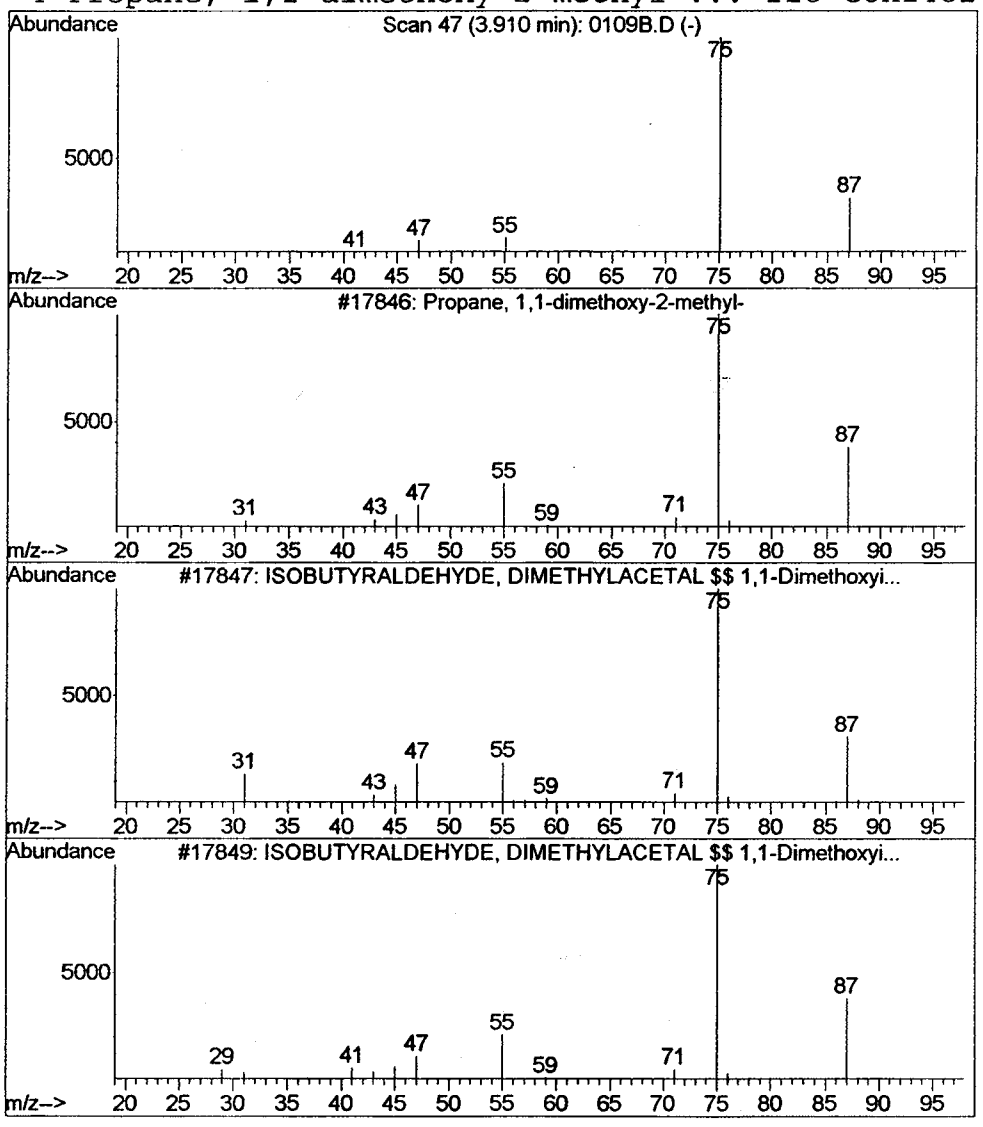
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

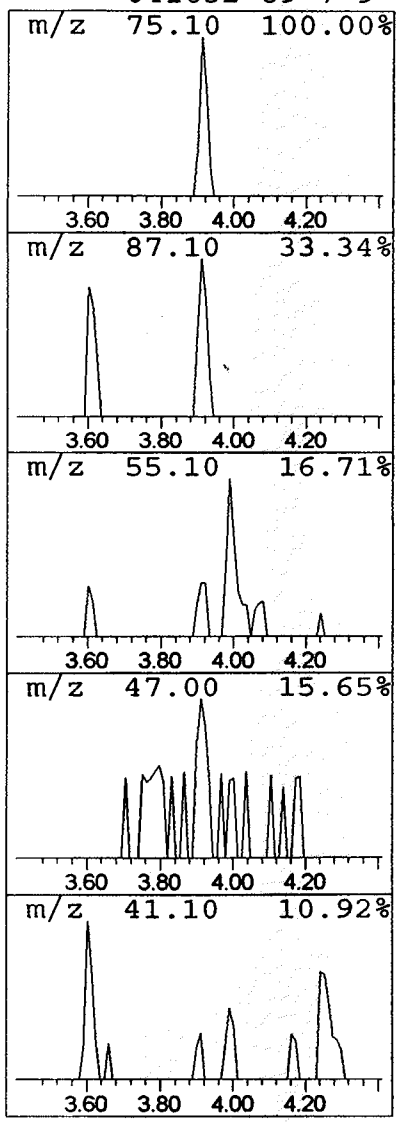
****

Peak Number 2 Propane, 1,1-dimethoxy-2-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.91	0.05 ug/L	21080	1,4-Dichlorobenzene-d4	9.71

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





NYC-WTC_000051518

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

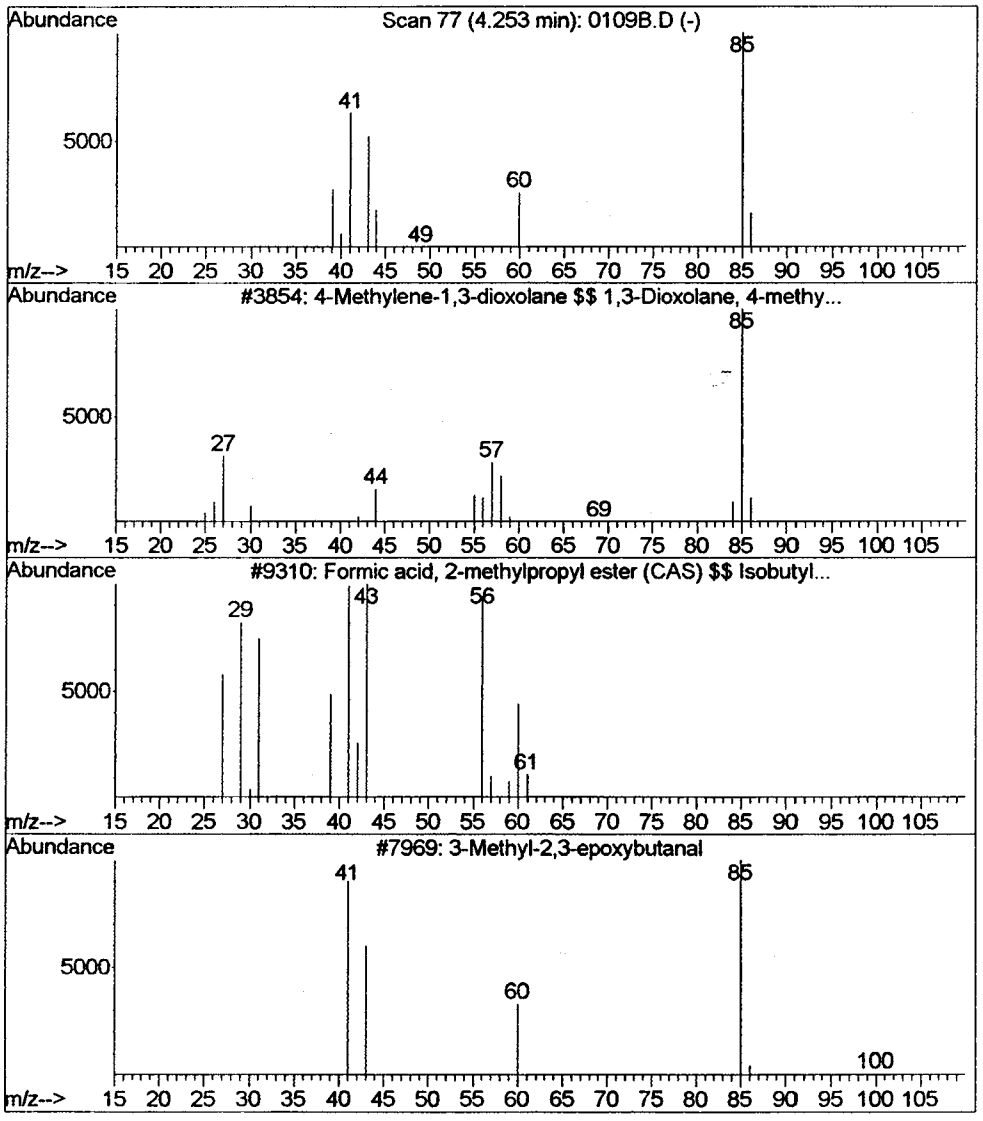
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

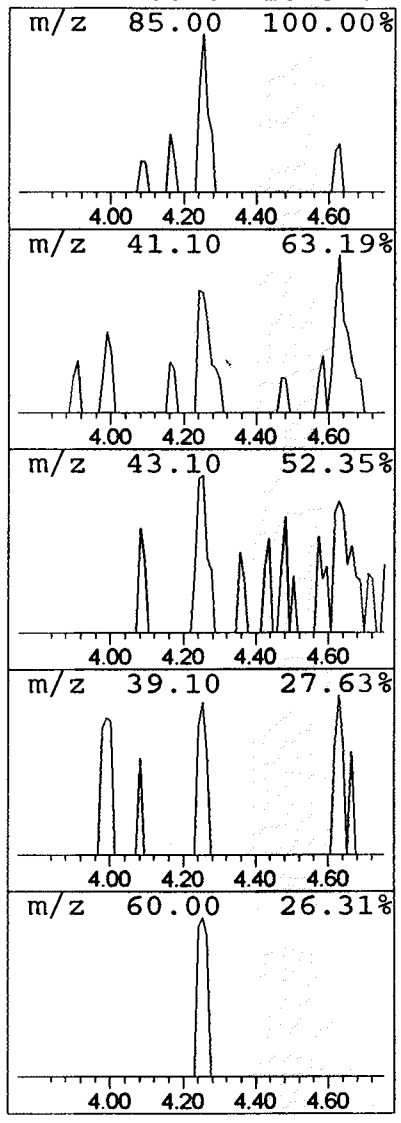
****

Peak Number 3 4-Methylene-1,3-dioxolane $... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.25	0.06 ug/L	28388	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methylene-1,3-dioxolane $$ 1,3...	86	C4H6O2	004362-24-7	9
2	Formic acid, 2-methylpropyl este...	102	C5H10O2	000542-55-2	9
3	3-Methyl-2,3-epoxybutanal	100	C5H8O2	000000-00-0	9
4	ETHYLENE DIAMMONIUM DICHLORIDE	60	C2H8N2	000107-15-3	5





NYC-WTC_000051519

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

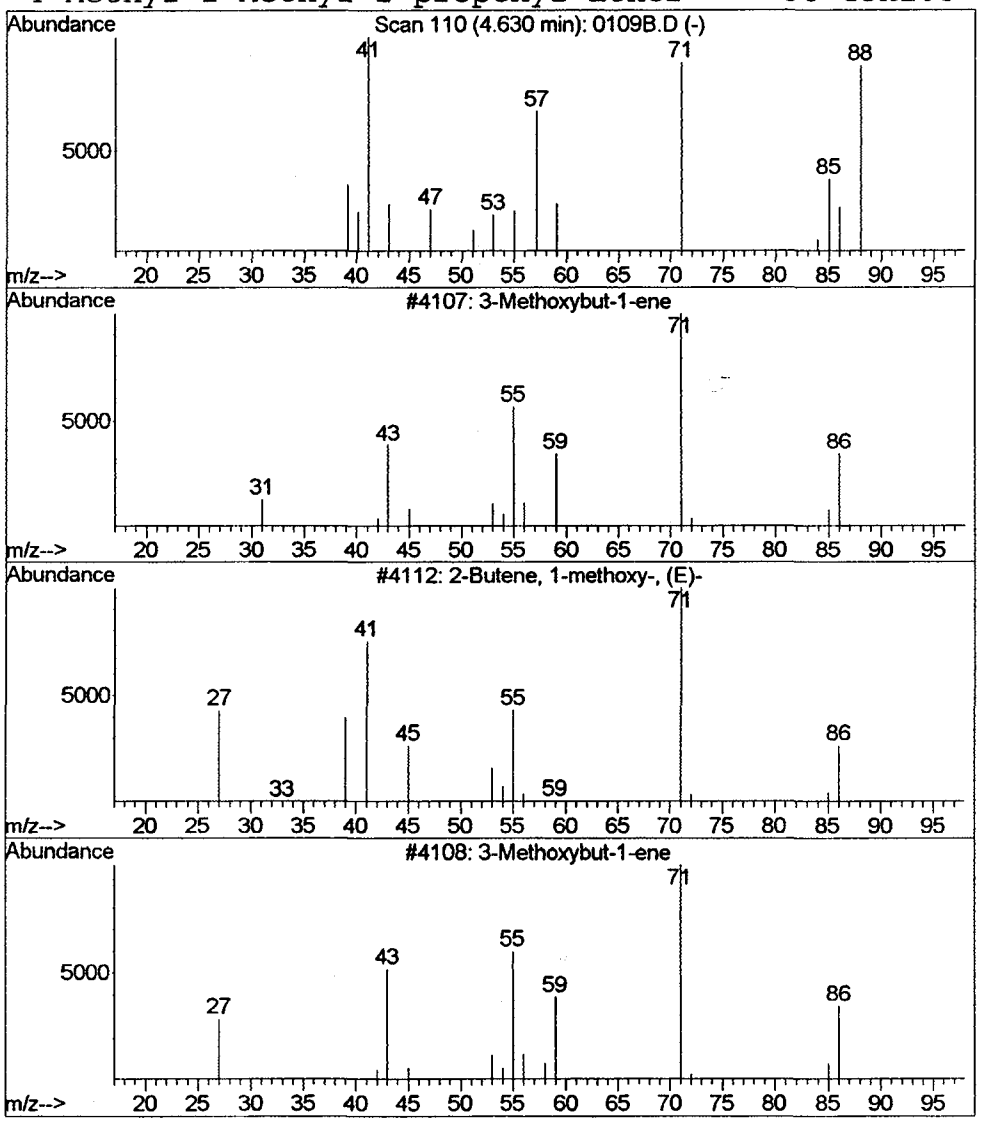
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

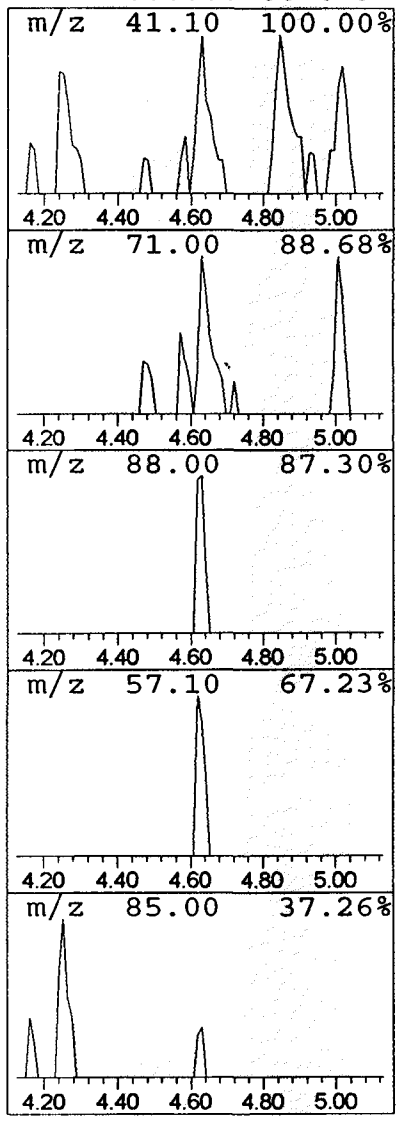
****

Peak Number 4 3-Methoxybut-1-ene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.63	0.08 ug/L	36317	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methoxybut-1-ene	86	C5H10O	017351-24-5	17
2	2-Butene, 1-methoxy-, (E)-	86	C5H10O	010034-14-7	17
3	3-Methoxybut-1-ene	86	C5H10O	017351-24-5	9
4	Methyl 1-Methyl-2-propenyl Ether	86	C5H10O	000000-00-0	9





NYC-WTC_000051520

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

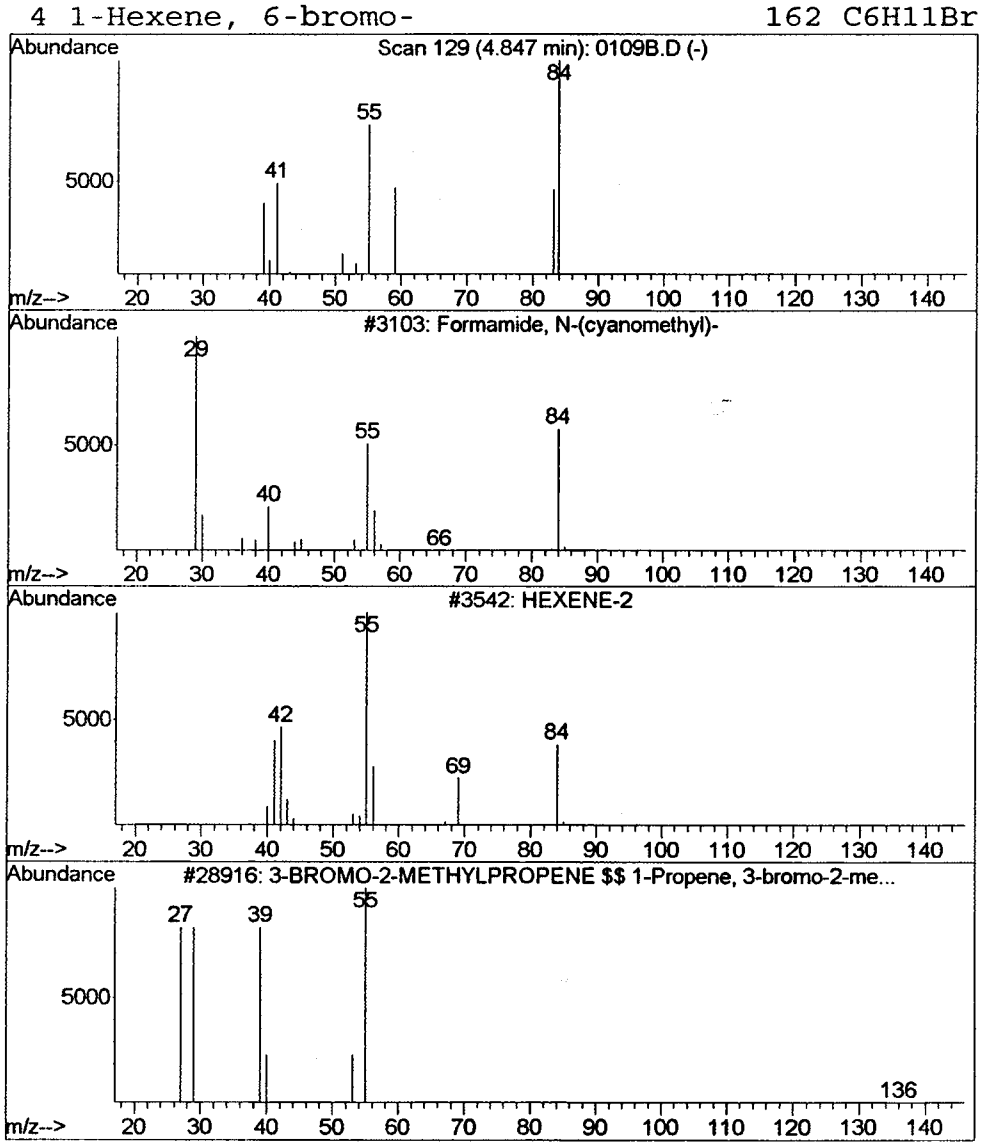
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

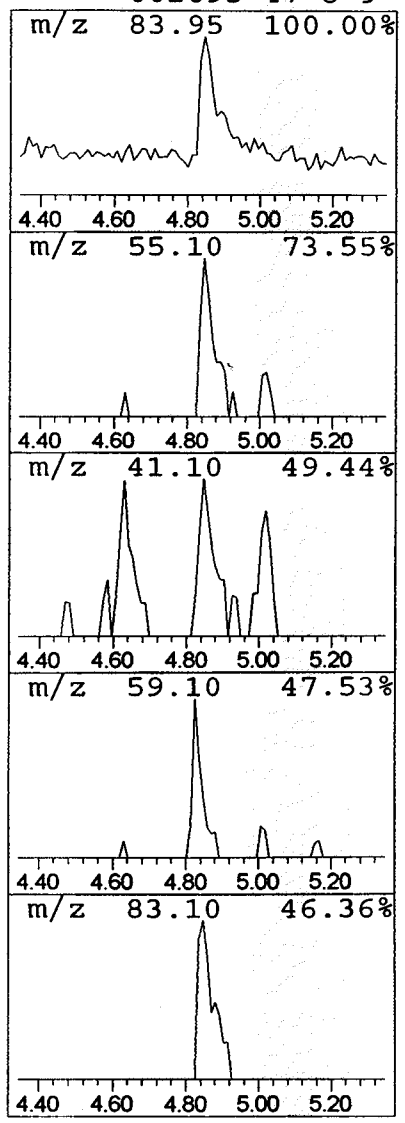
****

Peak Number 5 Formamide, N-(cyanomethyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.85	0.17 ug/L	76299	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Formamide, N-(cyanomethyl)-	84	C3H4N2O	005018-27-9	9
2	HEXENE-2	84	C6H12	000000-00-0	9
3	3-BROMO-2-METHYLPROPENE $$ 1-Pro...	134	C4H7Br	001458-98-6	9
4	1-Hexene, 6-bromo-	162	C6H11Br	002695-47-8	9





NYC-WTC_000051521

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

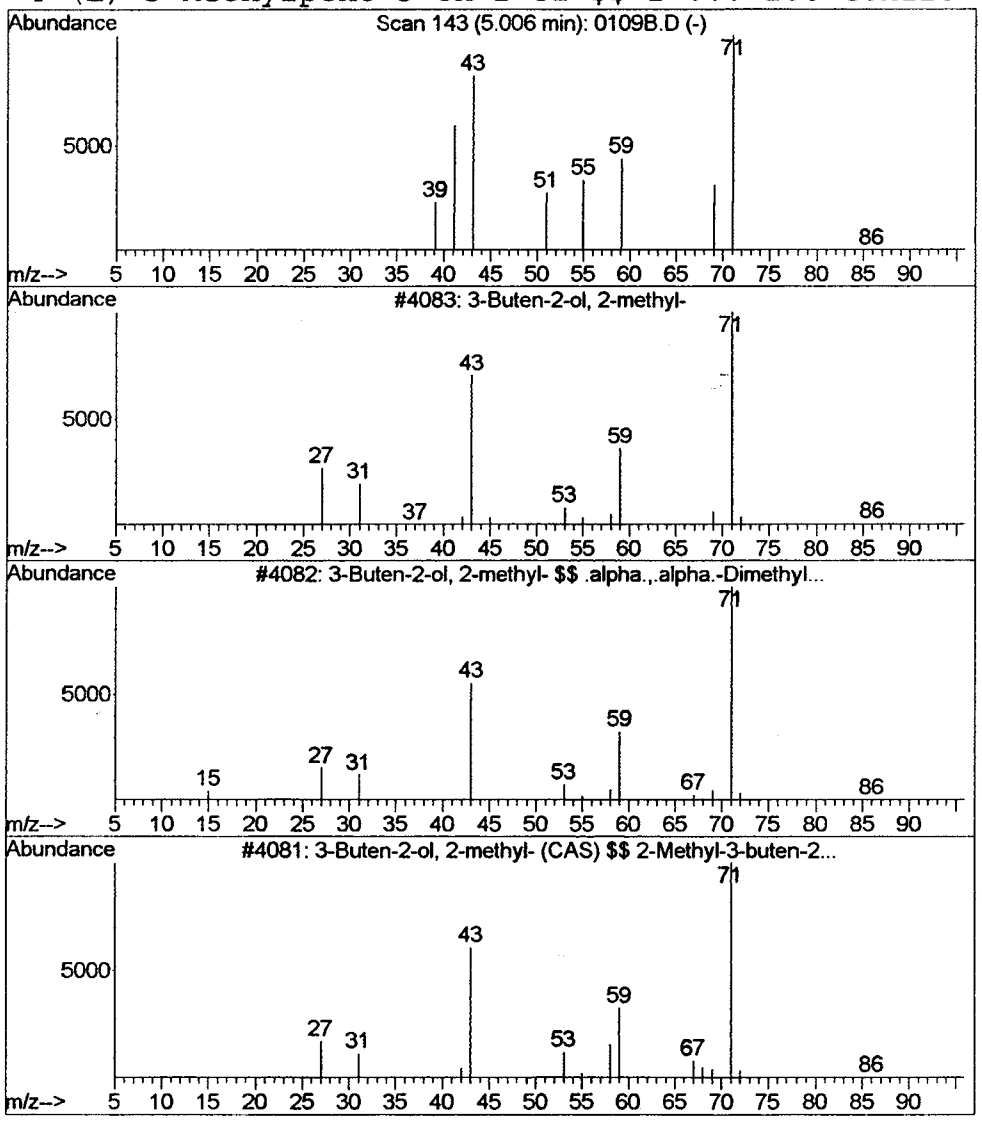
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

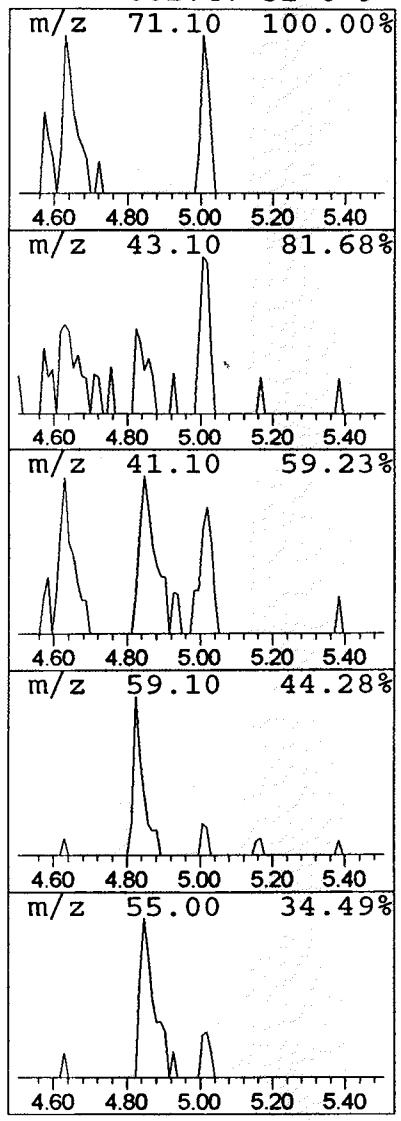
****

Peak Number 6 3-Buten-2-ol, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	0.06 ug/L	27816	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	9
2	3-Buten-2-ol, 2-methyl- $$ .alph...	86	C5H10O	000115-18-4	9
3	3-Buten-2-ol, 2-methyl- (CAS) $$...	86	C5H10O	000115-18-4	9
4	(E)-3-Methylpent-3-en-2-ol $$ 2-...	100	C6H12O	002747-52-6	9





NYC-WTC_000051522

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

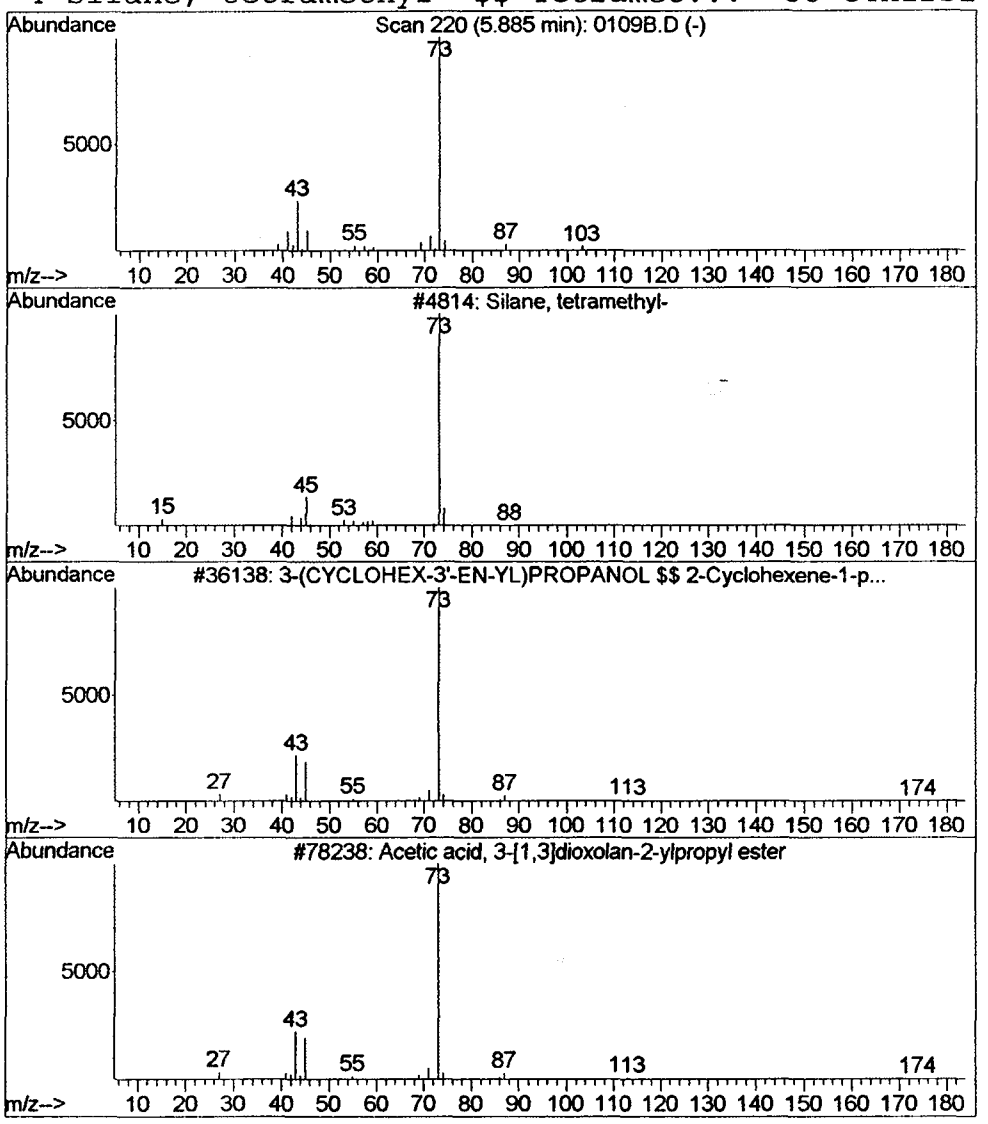
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

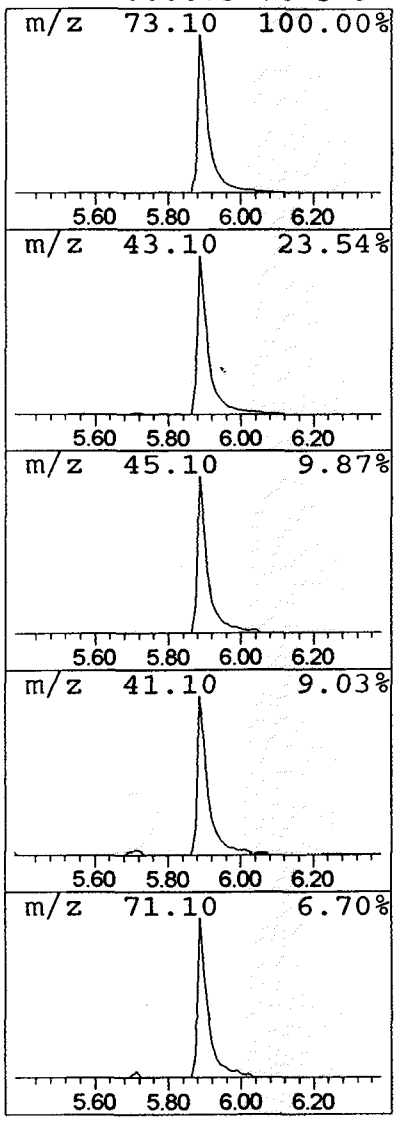
****

Peak Number 7 Silane, tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	3.77 ug/L	1645940	1,4-Dichlorobenzene-d4	9.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, tetramethyl-	88	C4H12Si	000075-76-3	37
2	3-(CYCLOHEX-3'-EN-YL)PROPANOL $$...	174	C8H14O4	015745-87-6	33
3	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	33
4	Silane, tetramethyl- $$ Tetramet...	88	C4H12Si	000075-76-3	9





NYC-WTC_000051523

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

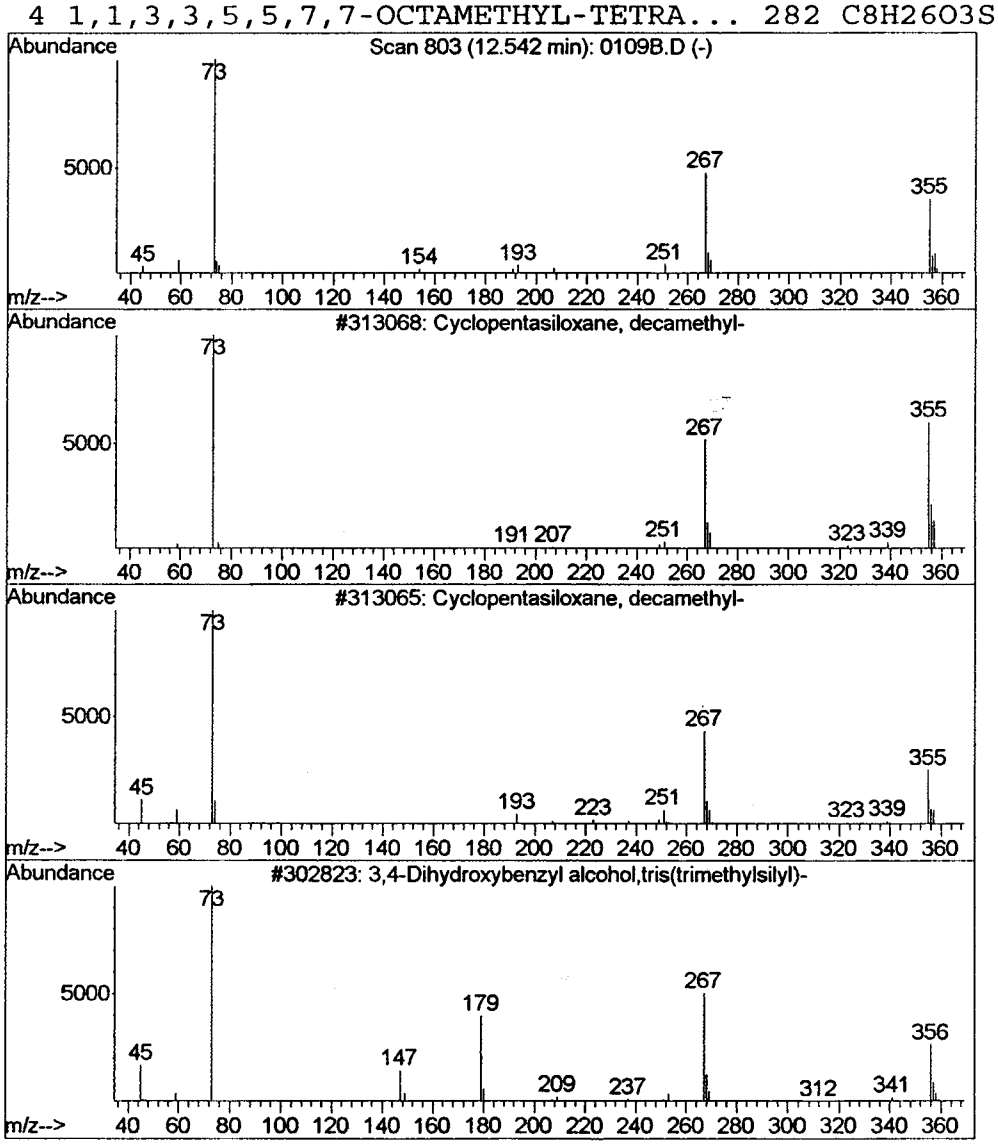
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

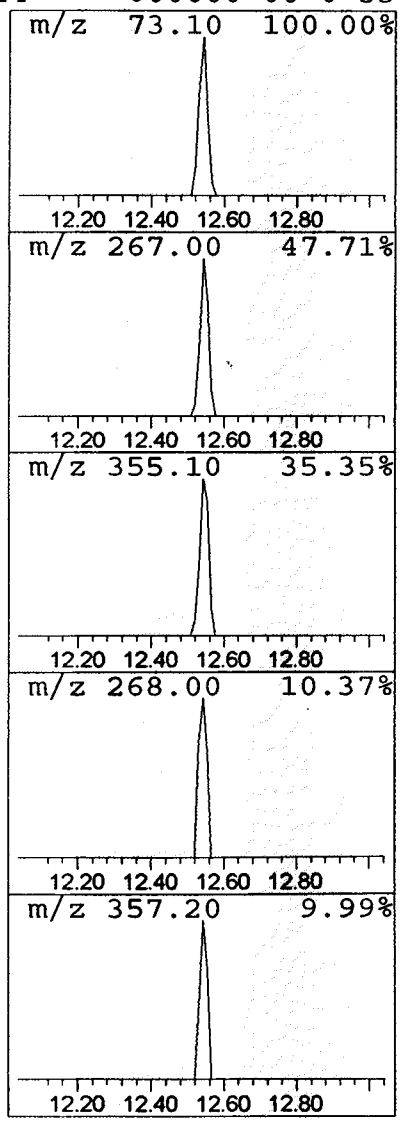
****

Peak Number 8 Cyclopentasiloxane, decamet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.13 ug/L	81429	Naphthalene-d8	13.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	83
2	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	64
3	3,4-Dihydroxybenzyl alcohol,tris...	356	C16H32O3Si3	068595-79-9	42
4	1,1,3,3,5,5,7,7-OCTAMETHYL-TETRA...	282	C8H26O3Si4	000000-00-0	33





NYC-WTC_000051524

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

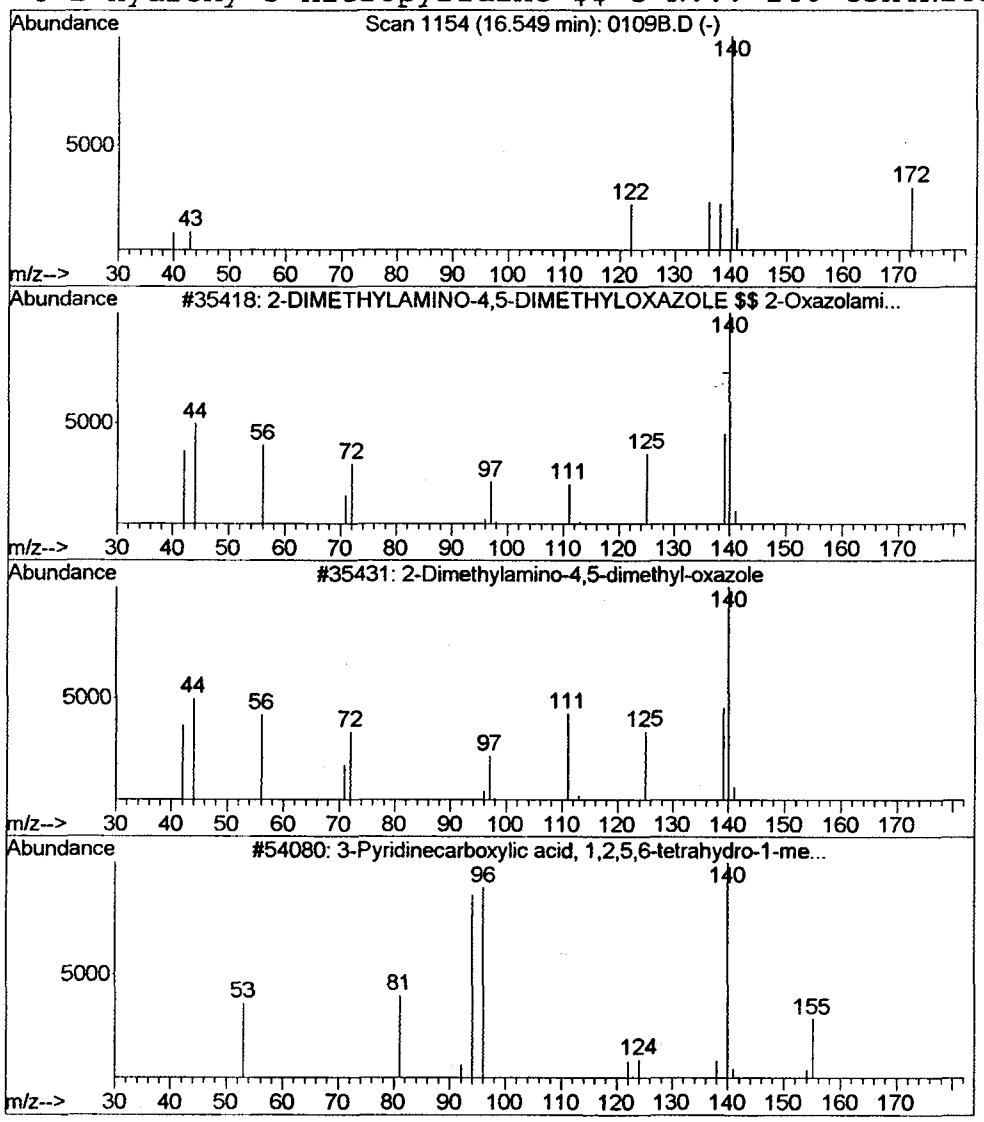
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

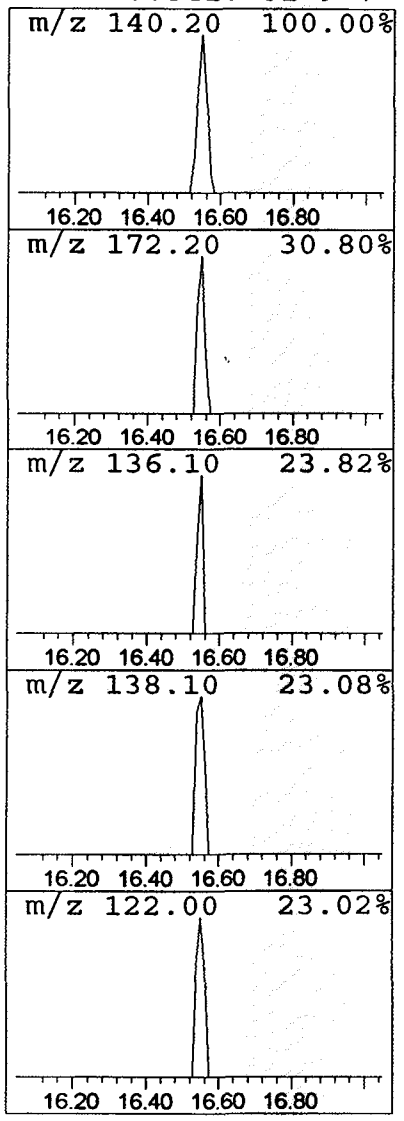
****

Peak Number 9 2-DIMETHYLAMINO-4,5-DIMETHY... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.55	0.03 ug/L	22910	Acenaphthene-d10	18.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-DIMETHYLAMINO-4,5-DIMETHYLOXAZ...	140	C7H12N2O	073777-29-4	9
2	2-Dimethylamino-4,5-dimethyl-oxa...	140	C7H12N2O	000000-00-0	9
3	3-Pyridinecarboxylic acid, 1,2,5...	155	C8H13NO2	000063-75-2	9
4	2-Hydroxy-5-nitropyridine $$ 5-N...	140	C5H4N2O3	005418-51-9	7





NYC-WTC_000051525

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

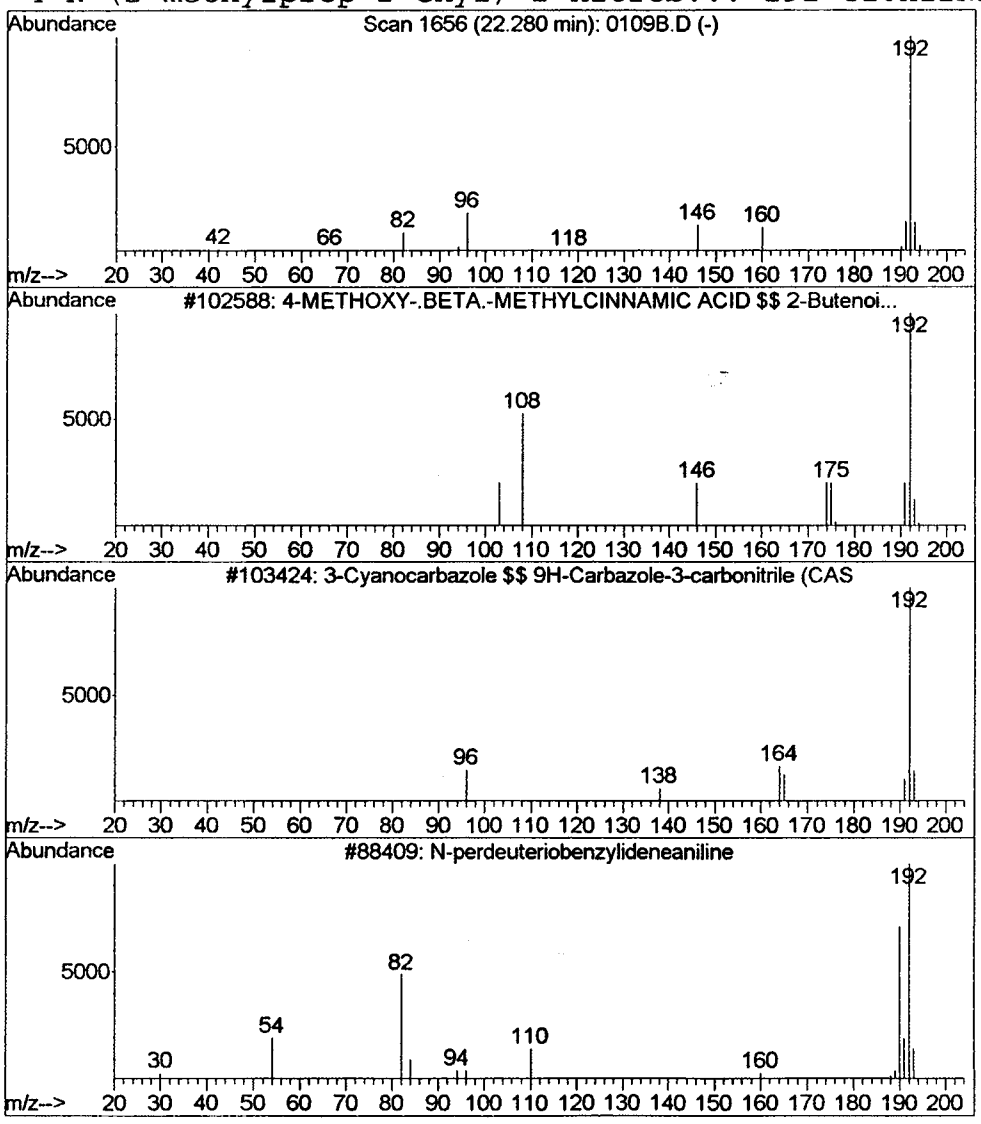
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

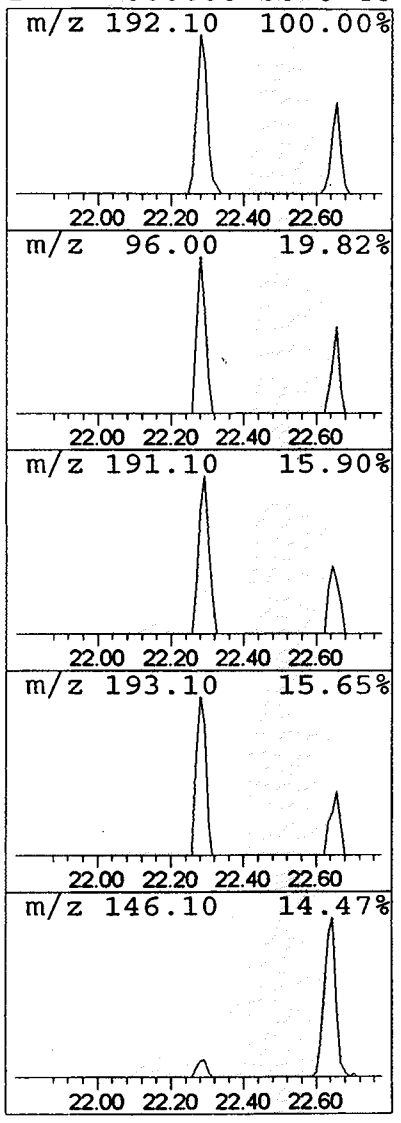
****

Peak Number 10 4-METHOXY-.BETA.-METHYLCINN... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.12 ug/L	89849	Phenanthrene-d10	22.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-METHOXY-.BETA.-METHYLCINNAMIC ...	192	C11H12O3	019169-94-9	72
2	3-Cyanocarbazole $$ 9H-Carbazole...	192	C13H8N2	057102-93-9	64
3	N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	58
4	N-(2-methylprop-2-enyl)-2-nitrob...	192	C10H12N2O2	000000-00-0	46





NYC-WTC_000051526

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

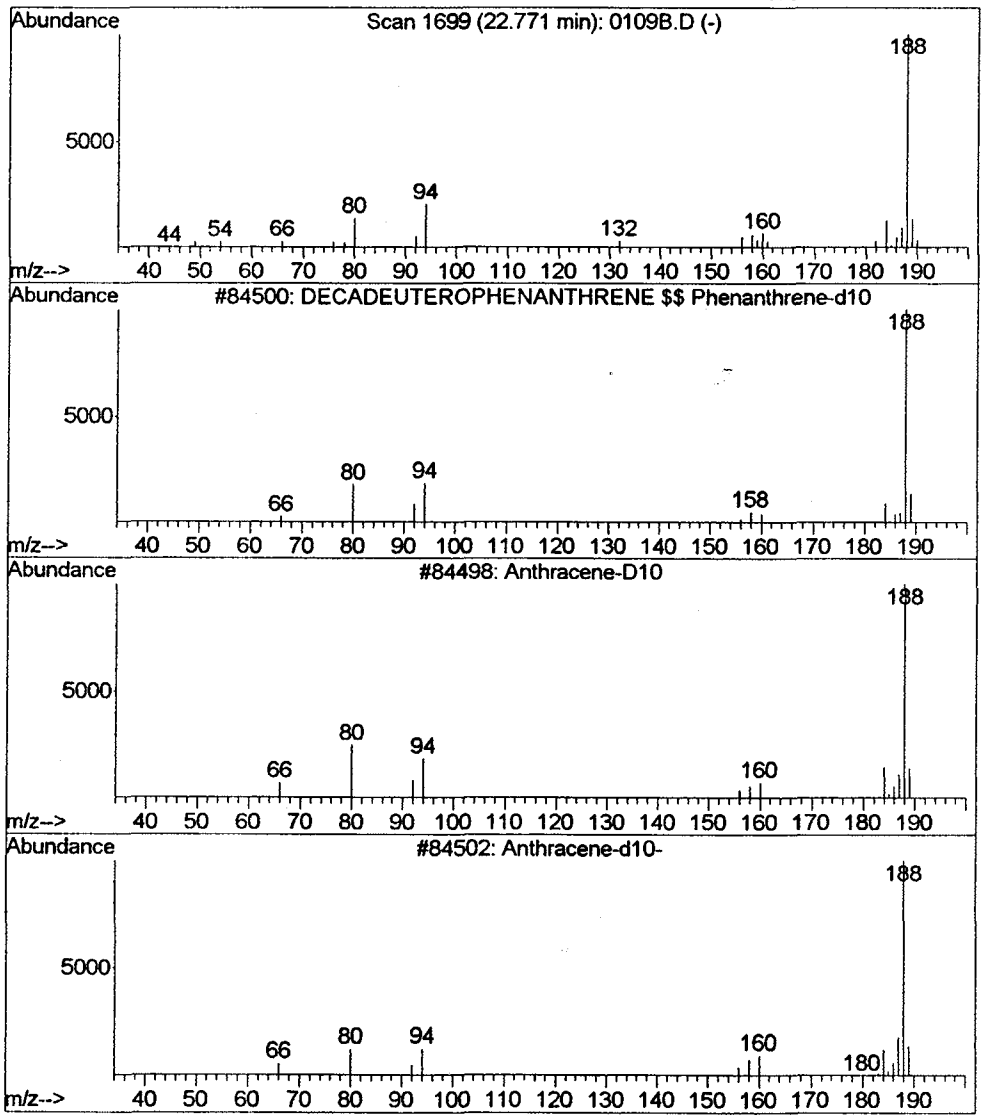
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

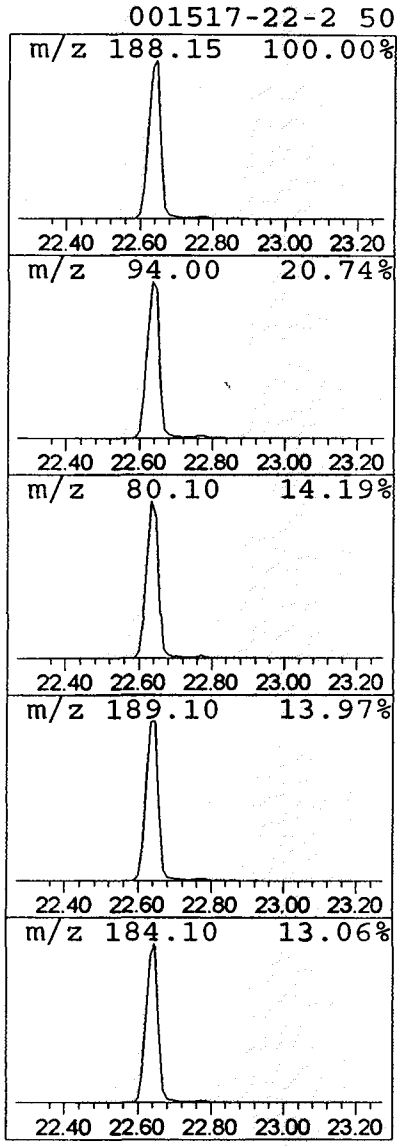
****

Peak Number 11 DECADEUTEROPHENANTHRENE $$ ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	0.29 ug/L	209940	Phenanthrene-d10	22.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	DECADEUTEROPHENANTHRENE $$ Phena...	188	C14D10	001517-22-2	87
2	Anthracene-D10	188	C14D10	000000-00-0	86
3	Anthracene-d10-	188	C14D10	001719-06-8	72
4	Phenanthrene-d10	188	C14D10	001517-22-2	50





NYC-WTC_000051527

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

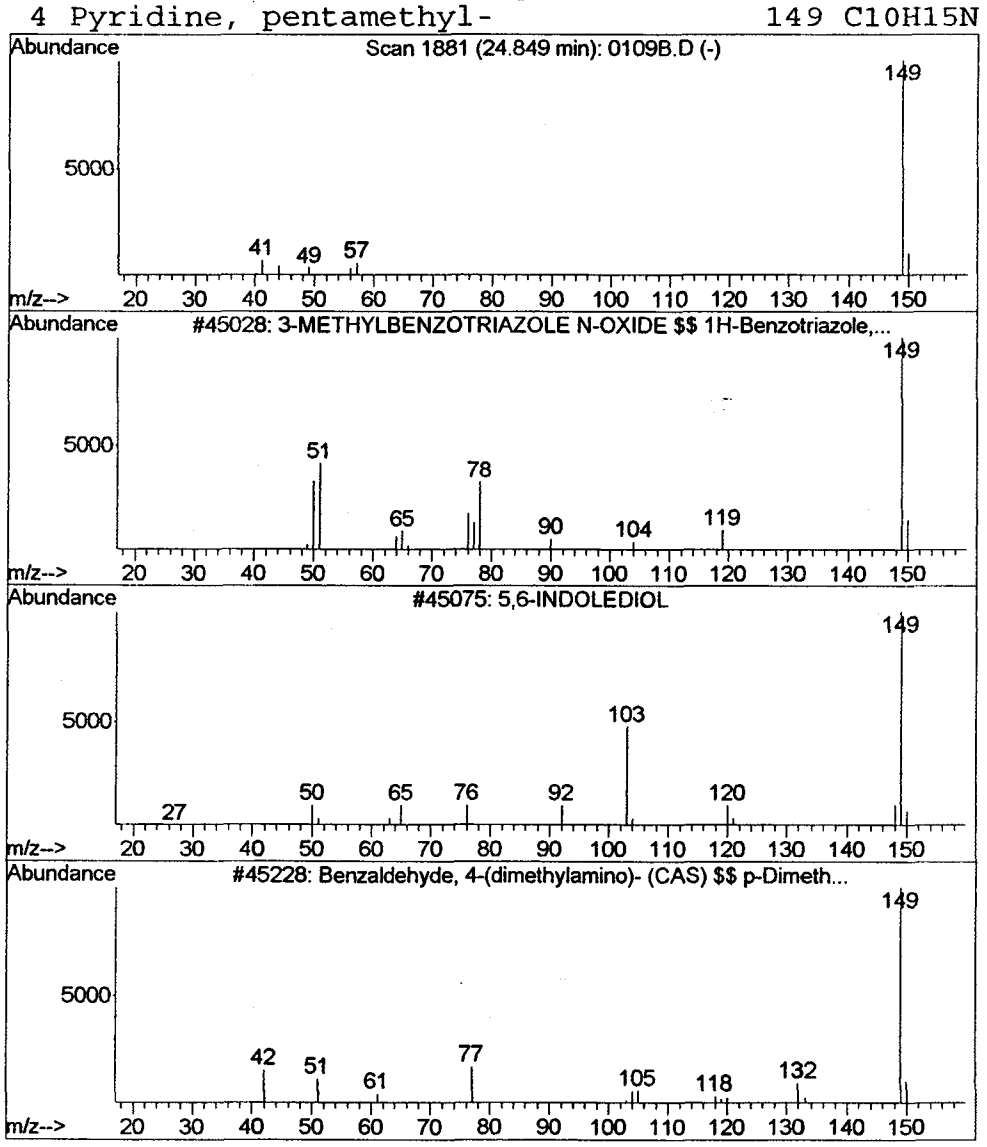
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

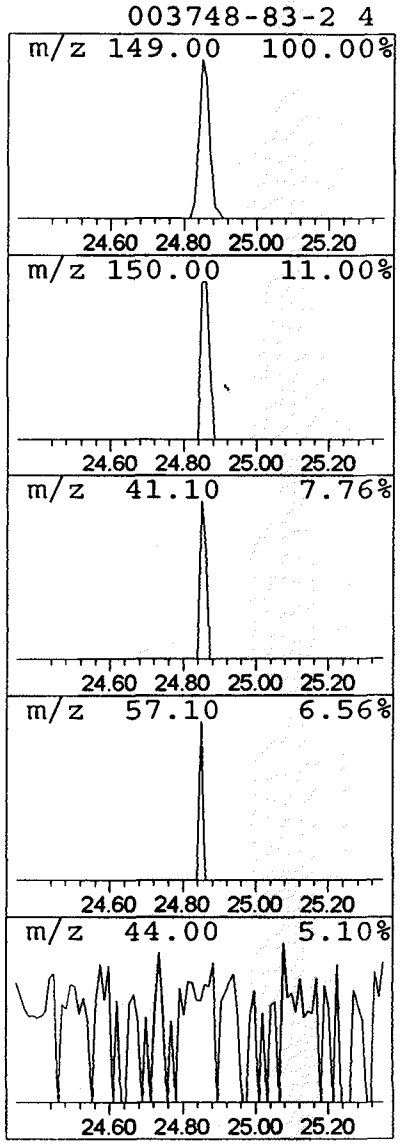
****

Peak Number 12 3-METHYLBENZOTRIAZOLE N-OXI... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.85	0.05 ug/L	35251	Phenanthrene-d10	22.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-METHYLBENZOTRIAZOLE N-OXIDE $$...	149	C7H7N3O	026198-26-5	9
2	5,6-INDOLEDIOL	149	C8H7NO2	000000-00-0	4
3	Benzaldehyde, 4-(dimethylamino)-...	149	C9H11NO	000100-10-7	4
4	Pyridine, pentamethyl-	149	C10H15N	003748-83-2	4





NYC-WTC_000051528

NYC 9/11 Public Portal Document

Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02JAN29\0109B.D Vial: 1
Acq On : 29 Jan 2002 10:31 am Operator:
Sample : lab blank 1/9 Inst : Instrumen
Misc : January 29, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

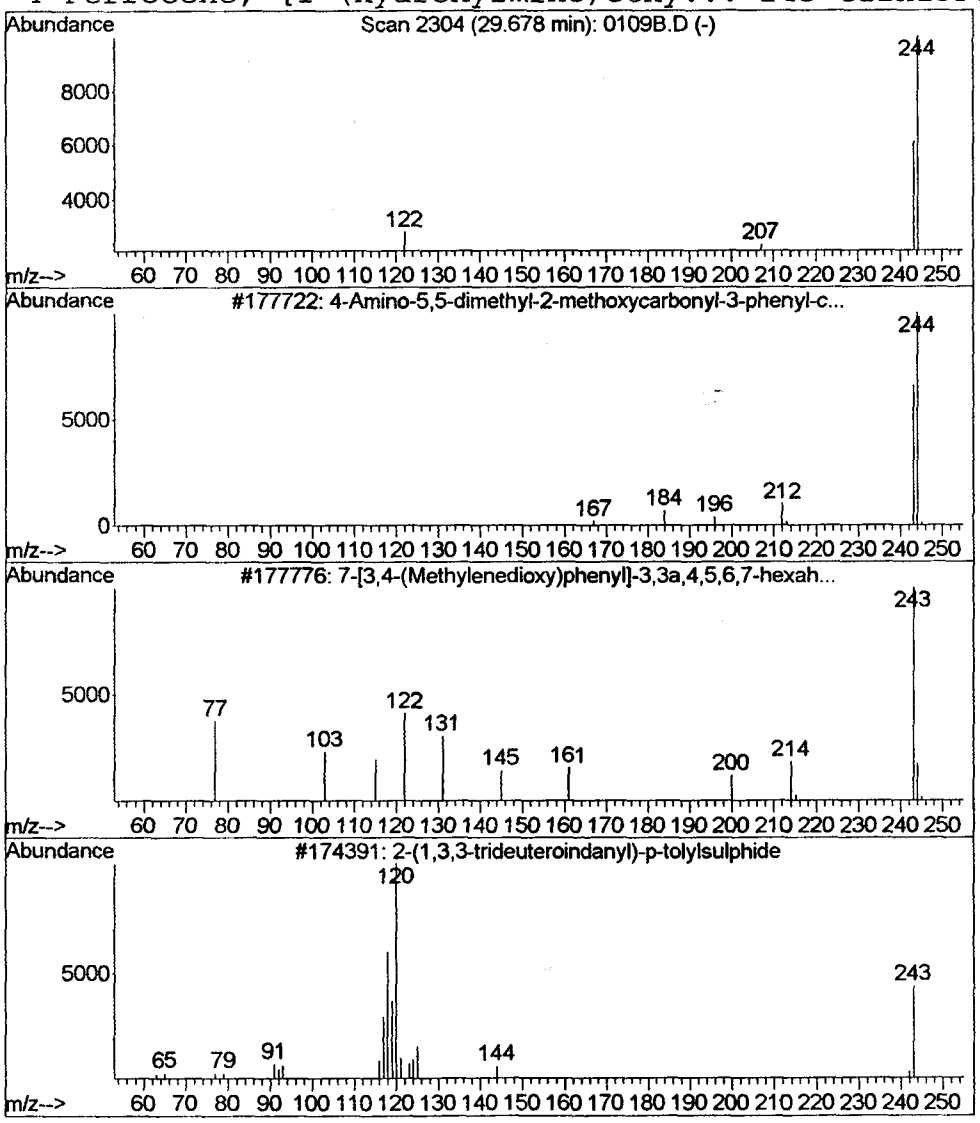
Quant Method : C:\MSDCHEM\1...\SCAN508.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

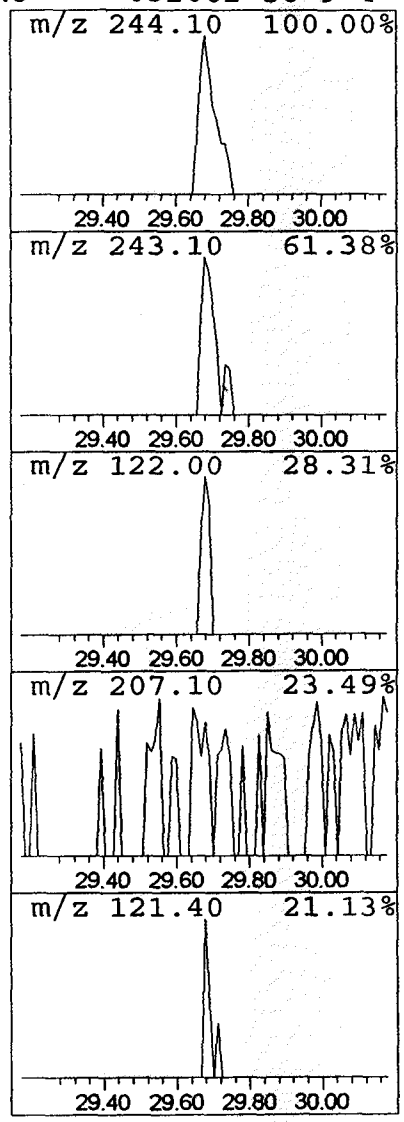
****

Peak Number 13 4-Amino-5,5-dimethyl-2-meth... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.68	0.04 ug/L	24567	Chrysene-d12	30.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Amino-5,5-dimethyl-2-methoxyca...	243	C15H17NO2	118089-29-5	5
2	7-[3,4-(Methylenedioxy)phenyl]-3...	243	C15H17NO2	000000-00-0	5
3	2-(1,3,3-trideuteroindanyl)-p-to...	243	C16H13D3S	000000-00-0	4
4	Ferrocene, [1-(hydroxyimino)ethy...	243	C12H13FeNO	032662-56-9	4





NYC-WTC_000051529

NYC 9/11 Public Portal Document

Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 29 Jan 2002 10:31 am
Data File: C:\MSDCHEM\1\5973N\02JAN29\0109B.D
Name: lab blank 1/9
Misc: January 29, 2002
Method: C:\MSDCHEM\1\METHODS\METHODS\SCAN508.M (RTE Integrator)
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...	3.60	0.1	ug/L	51376	1	9.71	8739380	20.0
Propane, 1,1-dime...	3.91	0.0	ug/L	21080	1	9.71	8739380	20.0
1-Methylene-1,3-d...	4.25	0.1	ug/L	28388	1	9.71	8739380	20.0
3-Methoxybut-1-ene	4.63	0.1	ug/L	36317	1	9.71	8739380	20.0
Formamide, N-(cya...	4.85	0.2	ug/L	76299	1	9.71	8739380	20.0
3-Buten-2-ol, 2-m...	5.01	0.1	ug/L	27816	1	9.71	8739380	20.0
Silane, tetramethyl-	5.89	3.8	ug/L	1645940	1	9.71	8739380	20.0
Cyclopentasiloxan...	12.54	0.1	ug/L	81429	2	13.19	12499900	20.0
2-DIMETHYLAMINO-4...	16.55	0.0	ug/L	22910	3	18.34	13567500	20.0
1-METHOXY-.BETA.-...	22.28	0.1	ug/L	89849	4	22.65	14625500	20.0
DECADEUTEROPHENAN...	22.77	0.3	ug/L	209940	4	22.65	14625500	20.0
3-METHYLBENZOTRIA...	24.85	0.0	ug/L	35251	4	22.65	14625500	20.0
1-Amino-5,5-dimet...	29.68	0.0	ug/L	24567	5	30.50	13557500	20.0

NYC-WTC_000051530