User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 03/25/2002 Time: 14:52:59

Sample: AE03265
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 3/25/02 14:44 Ended: 3/25/02 14:44 Analyst: MF
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

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Mar 25 08:56:34 2002 Quant Results File: 5080302.RES

Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Thu Mar 14 13:02:27 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.55	150	3871776	20.00	ug/L	0.00
10) Naphthalene-d8	13.04	136	10214141	20.00	ug/L	0.00
17) Acenaphthene-d10	18.15	164	5683233	20.00	ug/L	0.00
29) Phenanthrene-d10	22.51	188	8713889	20.00	ug/L	0.02
38) Chrysene-d12	30.34	240	7454504	20.00	ug/L	0.02
44) Perylene-d12	34.21	264	4229377	20.00	ug/L	0.02
System Monitoring Compounds						
37) 2,4-ddd surr	27.44	235	377183	3.63	ug/L	0.00
Spiked Amount 5.000			Recovery = 72.60%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	0.00	74	0	N.D.	d	
3) bis(2-chloroethyl) ether	0.00	93	0	N.D.		
4) 1,3 - Dichlorobenze	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-Chloropr	0.00	45	0	N.D.		
8) N-Nitroso-di-n-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) Naphthalene	0.00	128	0	N.D.		
16) Hexachlorobutadiene	0.00	225	0	N.D.		
18) Hexachlorocyclopentadiene	0.00	237	0	N.D.		
19) 2-chloronaphthalene	0.00	162	0	N.D.		
20) Dimethylphthalate	0.00	163	0	N.D.		
21) 2,6-Dinitrotoluene	0.00	165	0	N.D.	d	
22) Acenaphthylene	0.00	152	0	N.D.		
23) Acenaphthene	0.00	153	0	N.D.		
24) 2,4-Dinitrotoluene	0.00	165	0	N.D.		
25) Diethylphthalate	0.00	149	0	N.D.		
26) 4-Chlorophenyl-phenyl ethe	0.00	204	0	N.D.		
27) Fluorene	0.00	166	0	N.D.		
28) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.		
30) N-Nitrosodiphenylamine	0.00	169	0	N.D.		
31) 4-Bromophenyl-phenyl ether	0.00	248	0	N.D.		
32) Hexachlorobenzene	0.00	284	0	N.D.		
33) Phenanthrene	0.00	178	0	N.D.	d	
34) Anthracene	0.00	178	0	N.D.	d	
35) Di-n-butylphthalate	24.65	149	14858	0.03	ug/L	86
36) Fluoranthene	0.00	202	0	N.D.		
39) Pyrene	0.00	202	0	N.D.		
40) Butylbenzylphthalate	28.96	149	5599	0.03	ug/L #	16
41) Benzo (a) antracene	0.00	228	0	N.D.	d	
42) Bis (2-ethylhexyl) phthala	30.91	149	3550	0.01	ug/L	81
43) Chrysene	0.00	228	0	N.D.	d	
45) Di-n-Octylphthalate	32.79	149	3710	0.01	ug/L	1
46) Benzo (b) fluoranthene	0.00	252	0	N.D.		

6.91/5 = 138%

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

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Mar 25 08:56:34 2002 Quant Results File: 5080302.RES

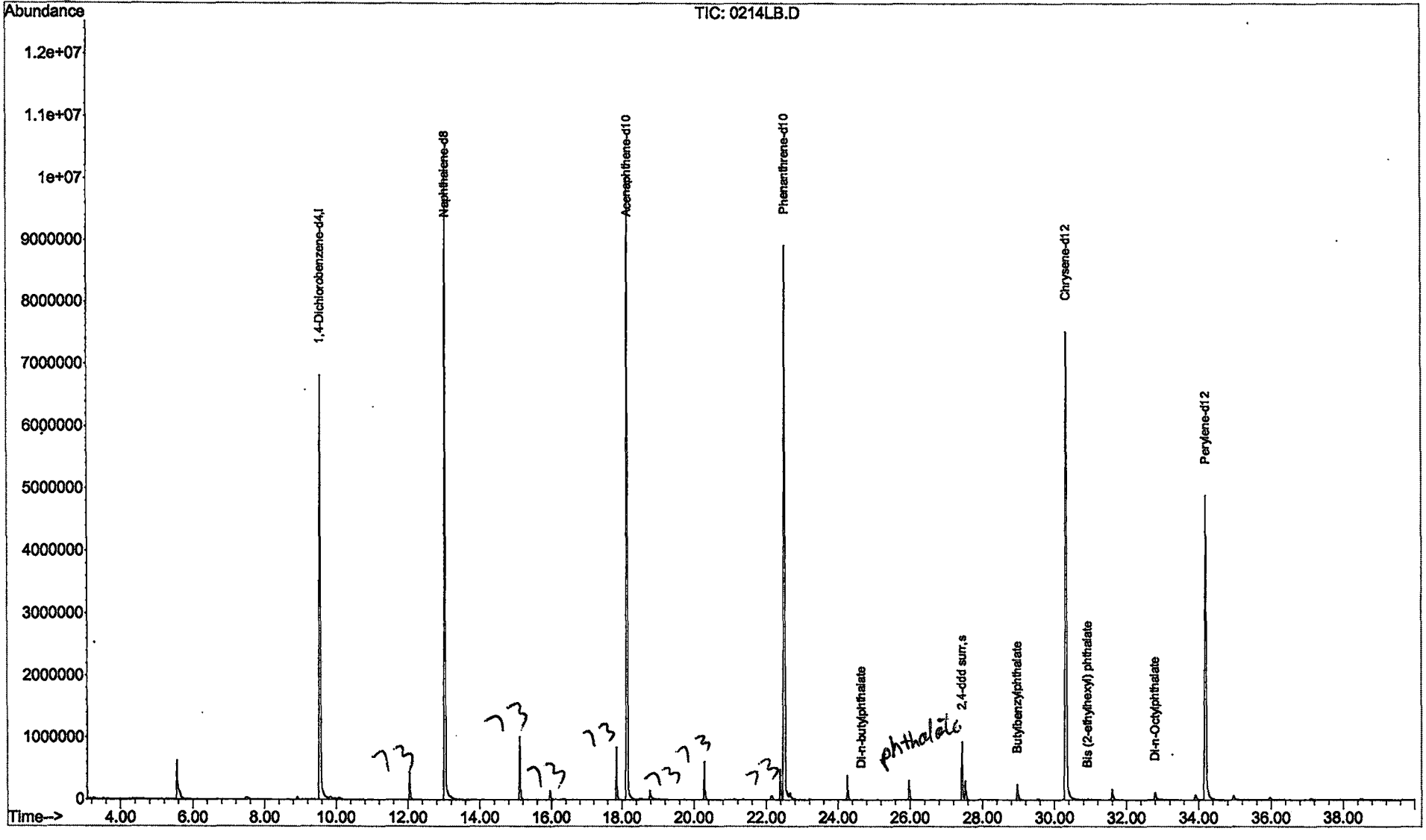
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Thu Mar 14 13:02:27 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
47) Benzo (k) fluoranthene	0.00	252	0	N.D.	
48) Benzo (a) pyrene	0.00	252	0	N.D.	d
49) Indeno (1,2,3-cd) pyrene	0.00	276	0	N.D.	
50) Dibenz (a,h) anthracene	0.00	278	0	N.D.	
51) Benzo (g,h,i) perylene	0.00	276	0	N.D.	

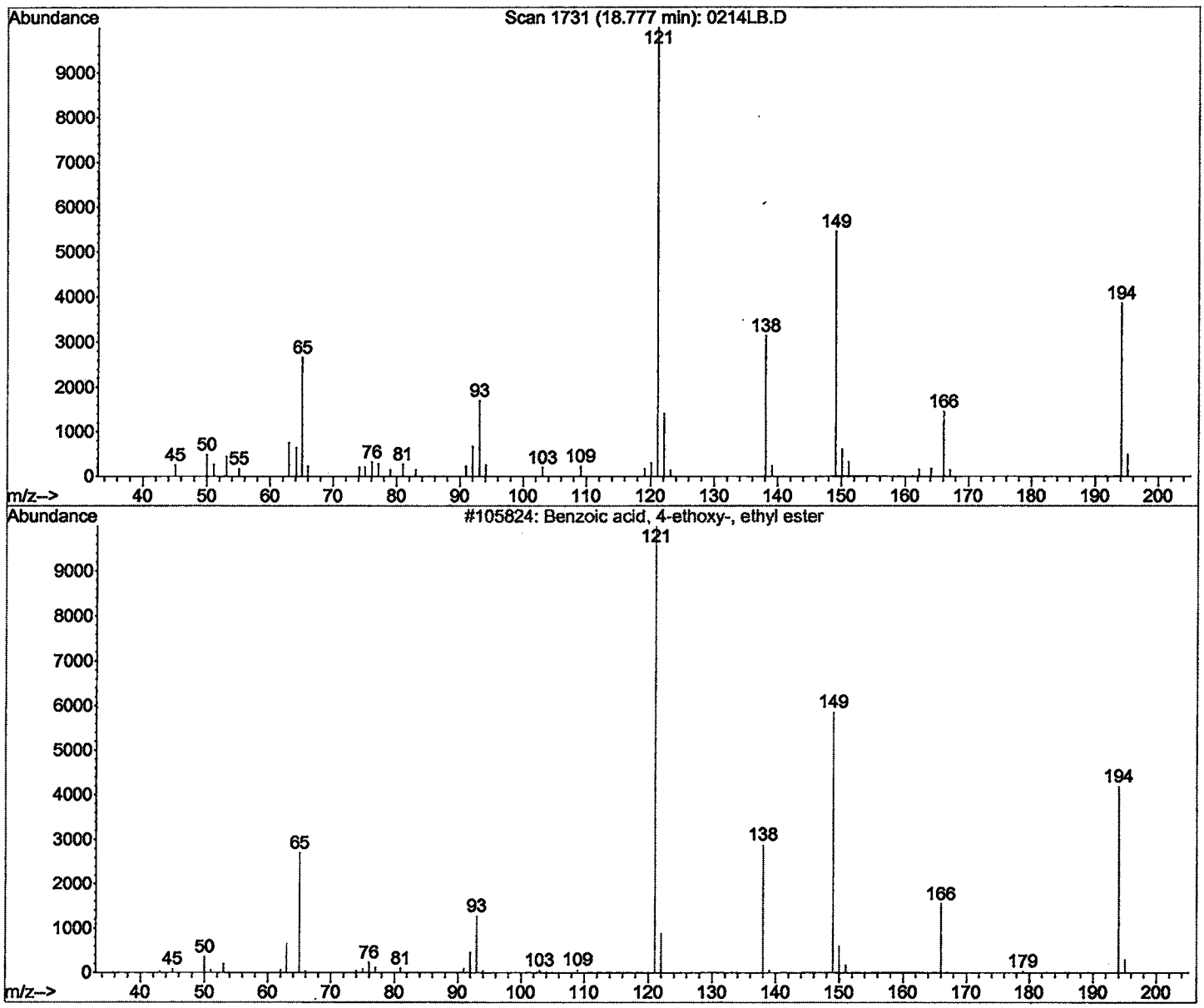
(#) = qualifier out of range (m) = manual integration
0214LB.D 5080302.M Mon Mar 25 08:57:30 2002 Page 2

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Mar 25 8:57 2002 Quant Results File: 5080302.RES

Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Thu Mar 14 13:02:27 2002
Response via : Initial Calibration



Library Searched : C:\Database\wiley7n.l
Quality : 97
ID : Benzoic acid, 4-ethoxy-, ethyl ester



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

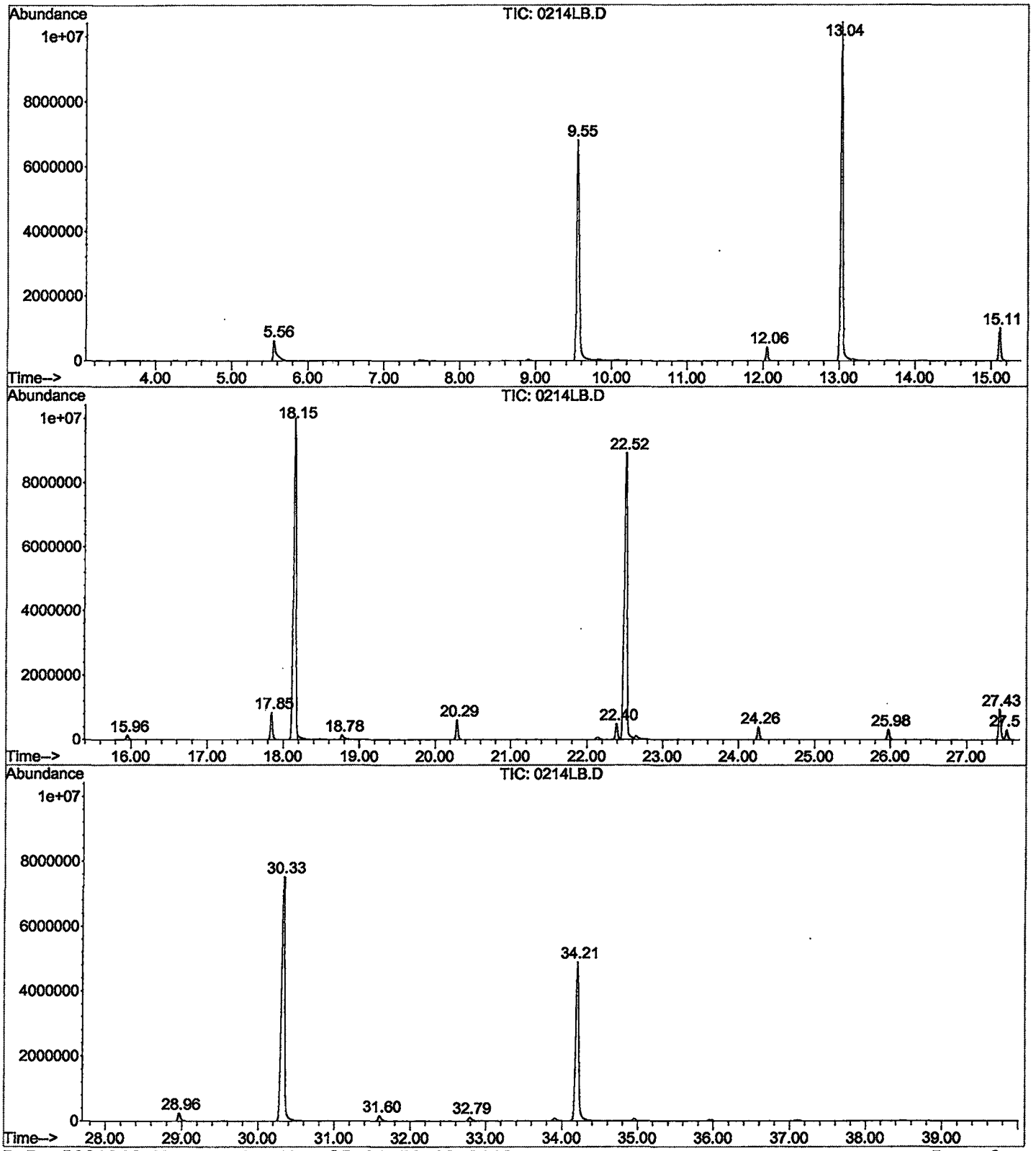
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.561	270	274	297	rBV	625371	1833317	7.64%	1.339%
2	9.552	708	714	734	rBV	6809543	16543136	68.90%	12.085%
3	12.056	985	990	1001	rBV	424759	748627	3.12%	0.547%
4	13.035	1091	1098	1112	rBV	10433717	22057719	91.87%	16.113%
5	15.112	1322	1327	1338	rBV	1011617	1849616	7.70%	1.351%
6	15.956	1416	1420	1425	rBV	145880	276849	1.15%	0.202%
7	17.852	1624	1629	1641	rBV	844863	1524384	6.35%	1.114%
8	18.151	1654	1662	1666	rBV	10012555	23282030	96.96%	17.008%
9	18.777	1727	1731	1746	rBV	145596	397209	1.65%	0.290%
10	20.291	1892	1898	1907	rBB	616032	1202494	5.01%	0.878%
11	22.396	2124	2130	2135	rBV2	486712	970014	4.04%	0.709%
12	22.523	2135	2144	2155	rBV2	8919099	24010848	100.00%	17.540%
13	24.264	2331	2336	2347	rBB	391998	727692	3.03%	0.532%
14	25.978	2519	2525	2534	rBB	314503	640154	2.67%	0.468%
15	27.430	2681	2685	2692	rBV	941098	1946593	8.11%	1.422%
16	27.529	2692	2696	2704	rVB2	307320	683674	2.85%	0.499%
17	28.963	2849	2854	2866	rBV2	247275	621384	2.59%	0.454%
18	30.332	2996	3005	3029	rBV2	7517817	22566553	93.98%	16.485%
19	31.602	3136	3145	3156	rBV2	170438	498705	2.08%	0.364%
20	32.790	3270	3276	3286	rBV2	118025	359727	1.50%	0.263%
21	34.205	3423	3432	3453	rBV2	4890203	14150678	58.93%	10.337%

Sum of corrected areas: 136891403

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\031902\0214LB.D
Operator : McLean
Acquired : 21 Mar 2002 5:58 am using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 2/14/02 RE-INJ
Misc Info :
Vial Number: 33
Quant File :5080302.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

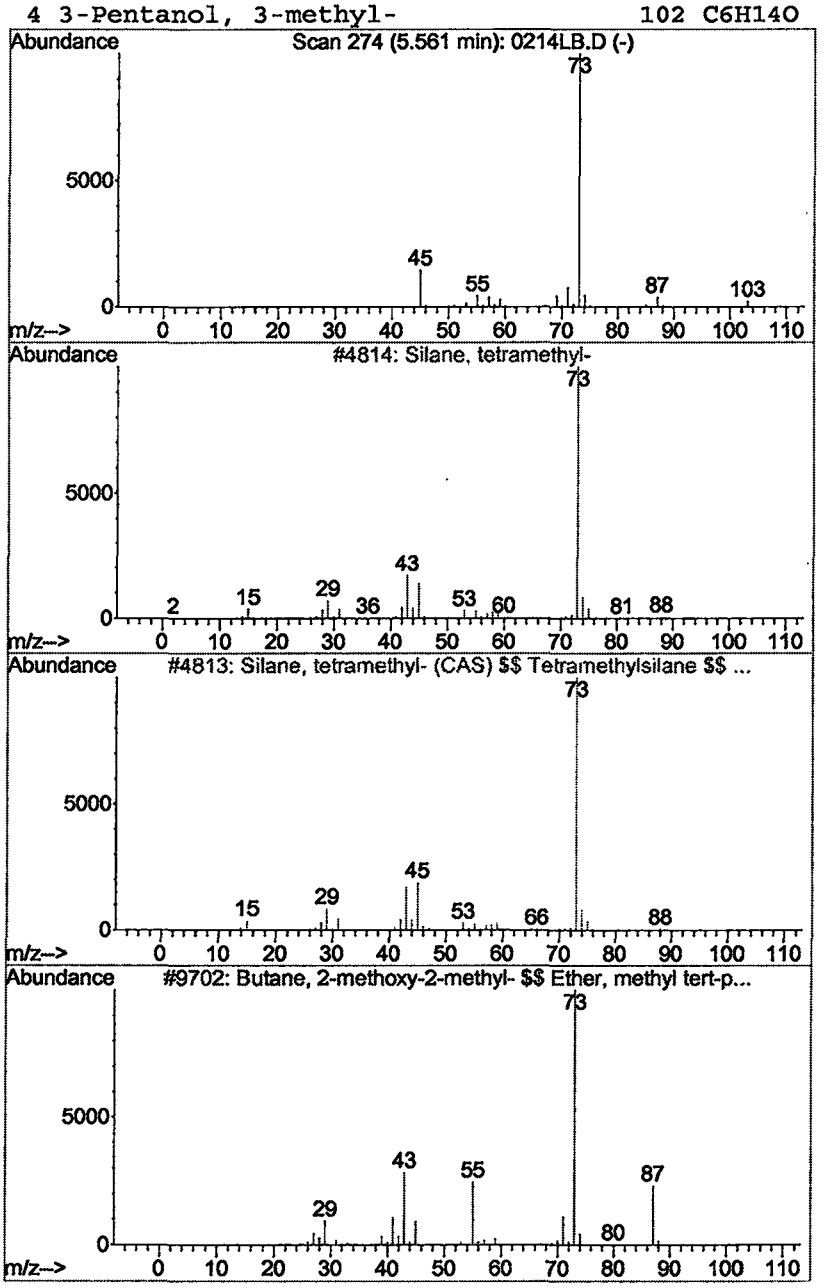
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

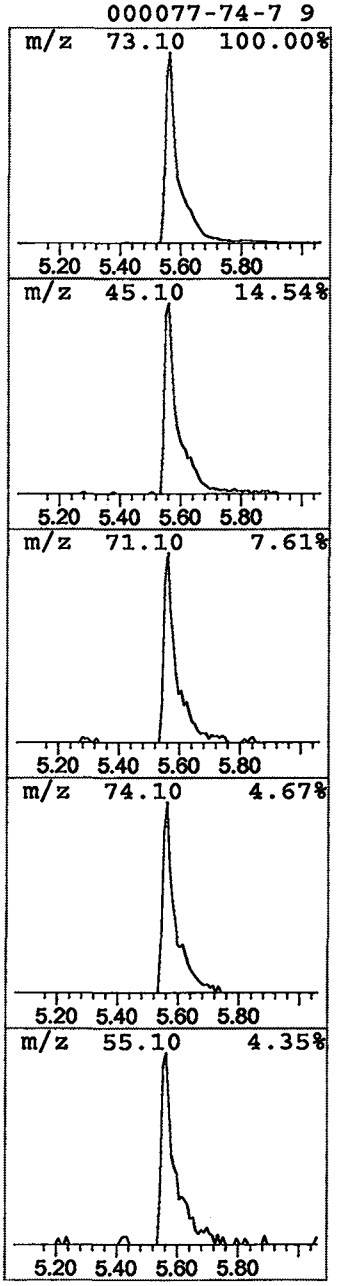
****

Peak Number 1 Silane, tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.56	2.22 ug/L	1833320	1,4-Dichlorobenzene-d4	9.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, tetramethyl-	88	C4H12Si	000075-76-3	47
2	Silane, tetramethyl- (CAS) $$ Te...	88	C4H12Si	000075-76-3	9
3	Butane, 2-methoxy-2-methyl- $$ E...	102	C6H14O	000994-05-8	9
4	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	9





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
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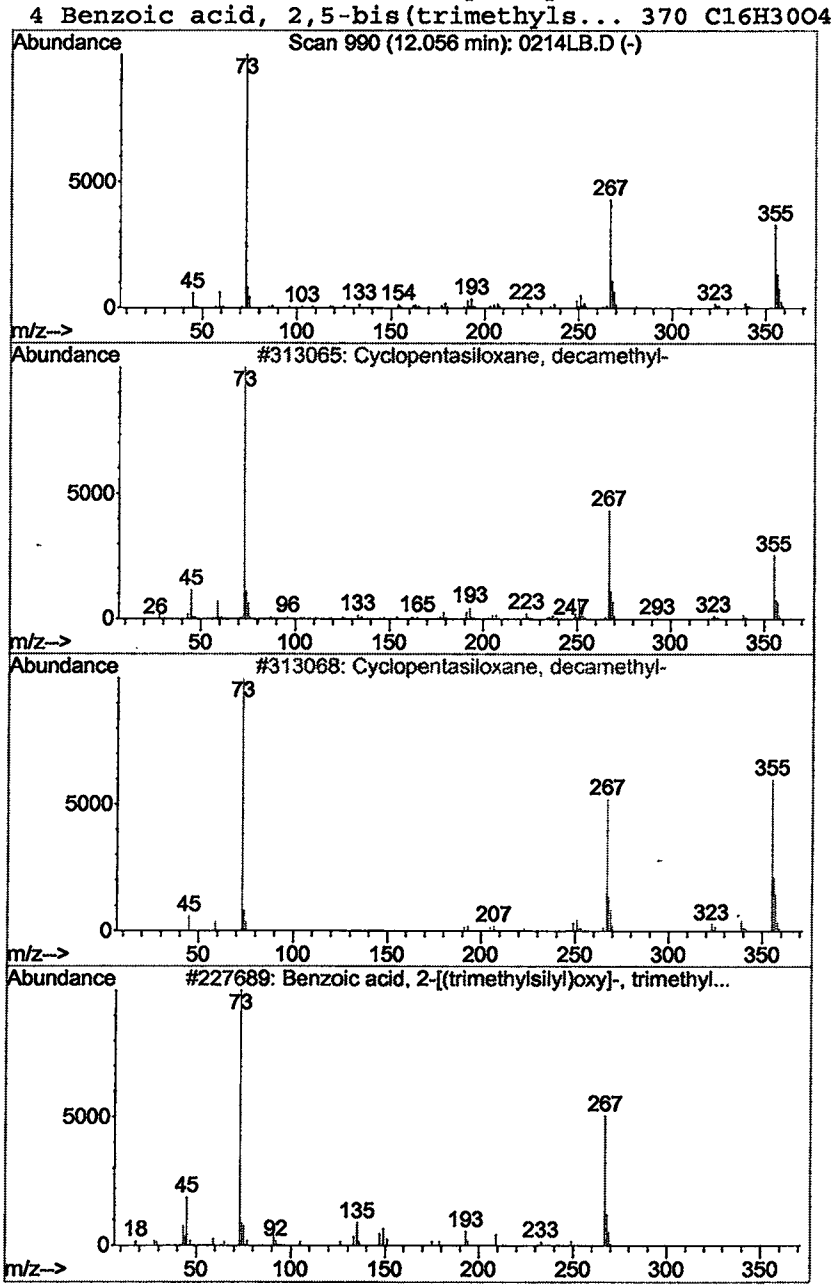
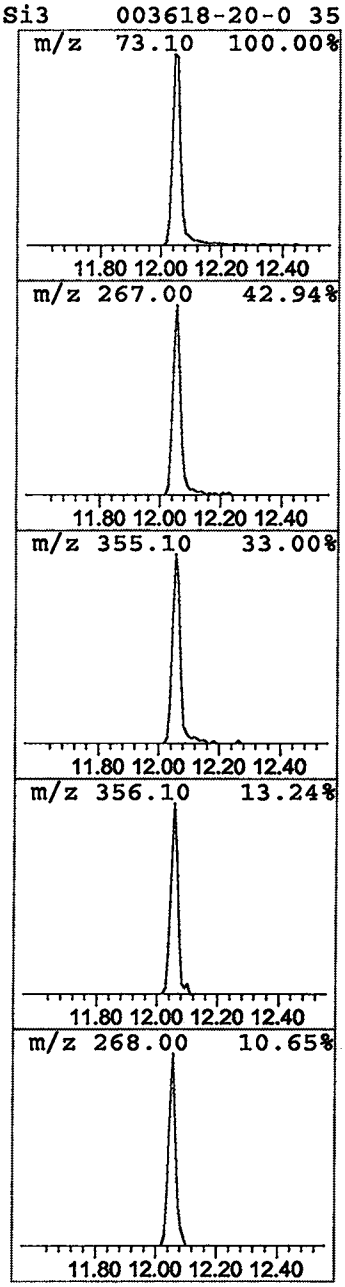
****

Peak Number 2 Cyclopentasiloxane, decamet... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.06	0.68 ug/L	748627	Naphthalene-d8	13.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
2	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	58
3	Benzoic acid, 2-[(trimethylsilyl...	282	C13H22O3Si2	003789-85-3	38
4	Benzoic acid, 2,5-bis(trimethyls...	370	C16H30O4Si3	003618-20-0	35





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

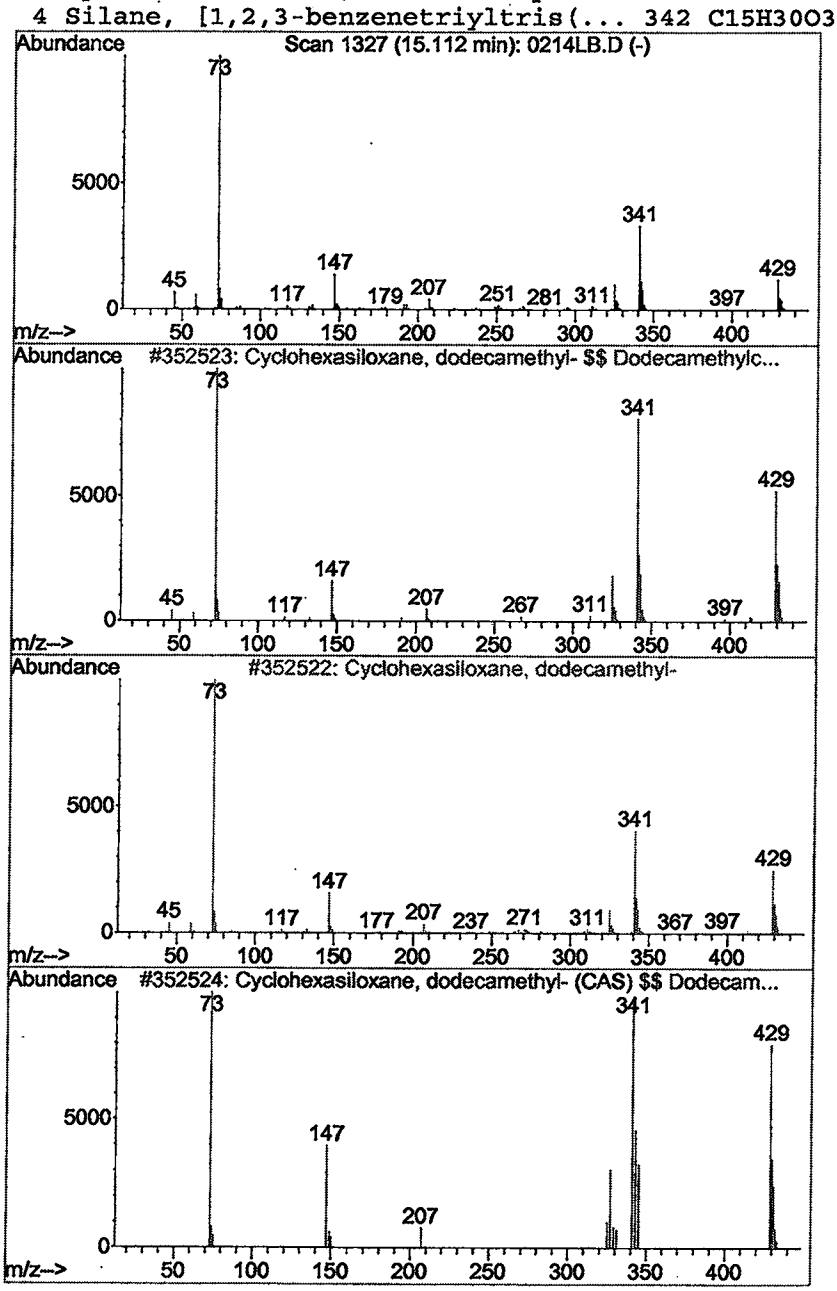
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

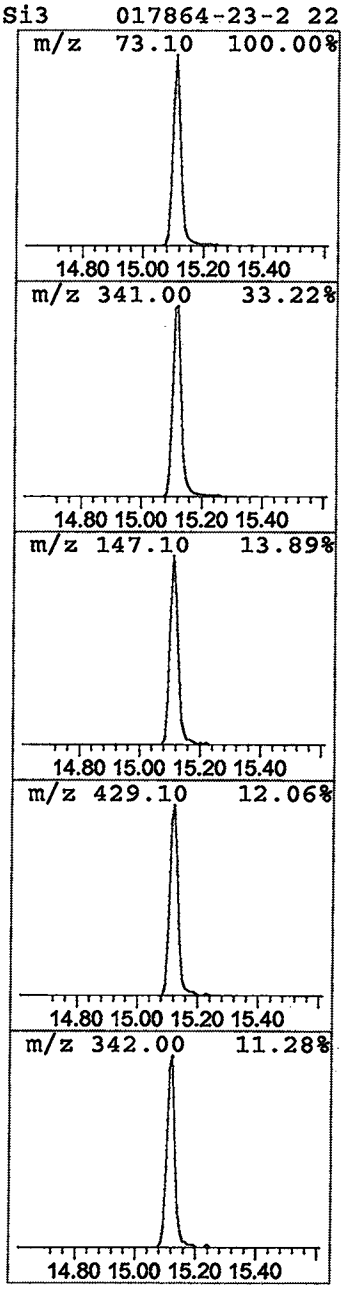
****

Peak Number 3 Cyclohexasiloxane, dodecame... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.11	1.68 ug/L	1849620	Naphthalene-d8	13.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexasiloxane, dodecamethyl-...	444	C12H36O6Si6	000540-97-6	91
2	Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	91
3	Cyclohexasiloxane, dodecamethyl-...	444	C12H36O6Si6	000540-97-6	23
4	Silane, [1,2,3-benzenetriyltris(...	342	C15H30O3Si3	017864-23-2	22





Library Search Compound Report

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Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

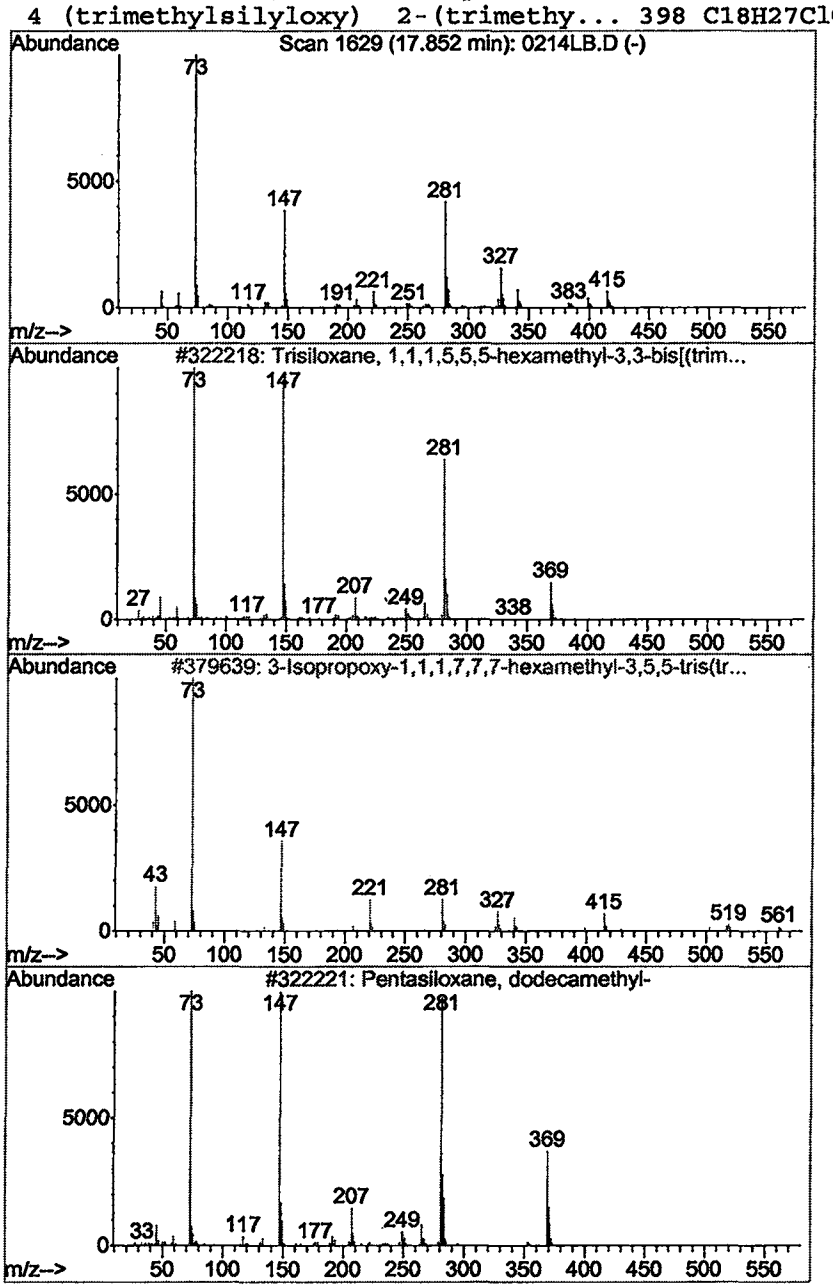
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

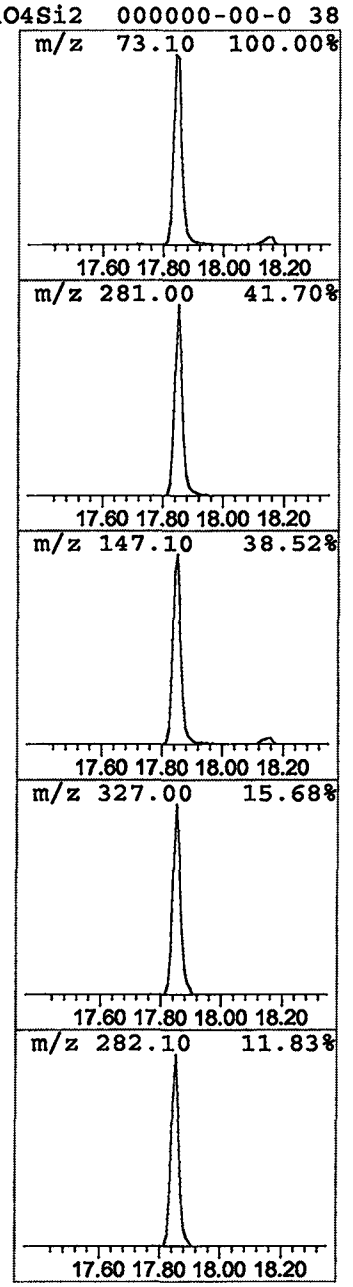
****

Peak Number 4 Trisiloxane, 1,1,1,5,5,5-he... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.85	1.31 ug/L	1524380	Acenaphthene-d10	18.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	40
2	3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	40
3	Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	39
4	(trimethylsilyloxy) 2-(trimethy...	398	C18H27ClO4Si2	000000-00-0	38





Library Search Compound Report

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Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

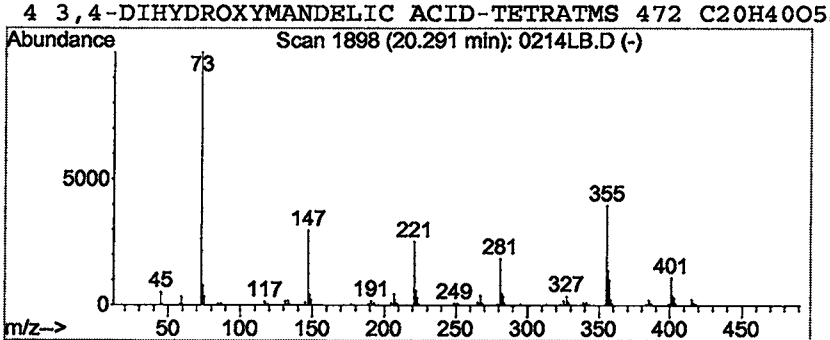
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

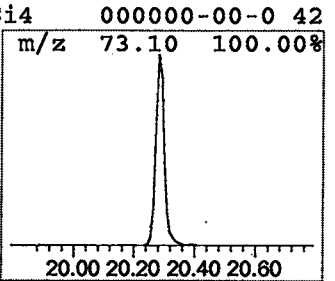
****

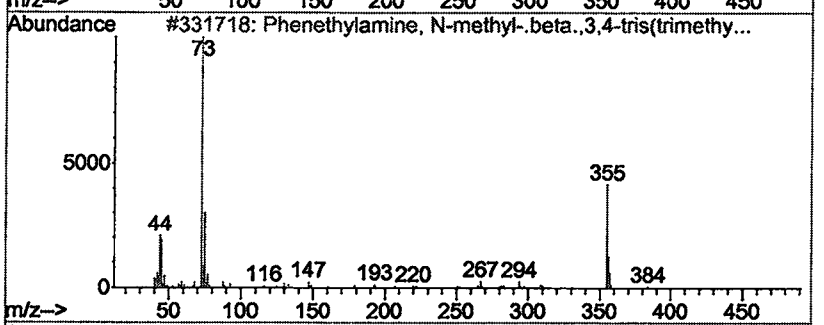
Peak Number 5 Phenethylamine, N-methyl-.b... Concentration Rank 4

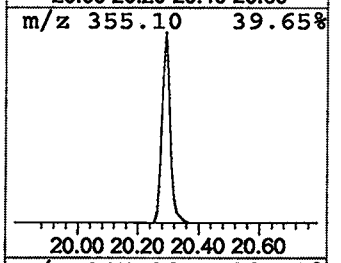
R.T.	EstConc	Area	Relative to ISTD	R.T.
20.29	1.03 ug/L	1202490	Acenaphthene-d10	18.15

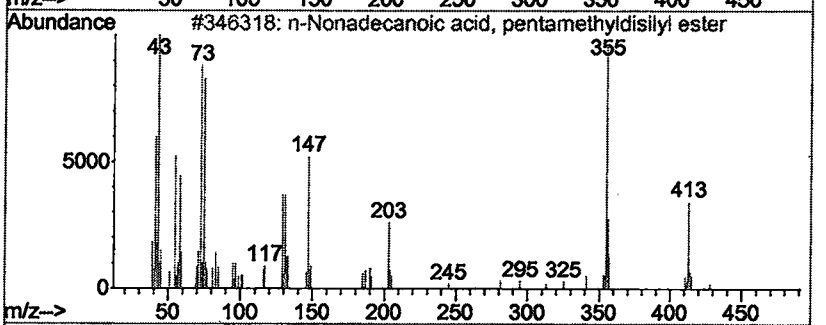
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenethylamine, N-methyl-.beta.,...	399	C18H37NO3Si3	010538-85-9	43
2	n-Nonadecanoic acid, pentamethyl...	428	C24H52O2Si2	000000-00-0	43
3	Benzeneacetic acid, .alpha.,3,4-...	472	C20H40O5Si4	037148-65-5	43
4	3,4-DIHYDROXYMANDELIC ACID-TETRATMS	472	C20H40O5Si4	000000-00-0	42

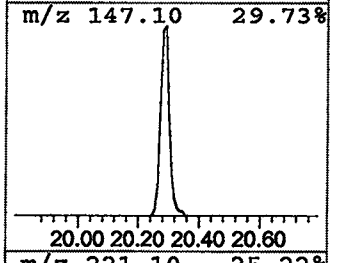


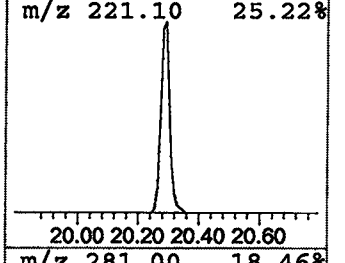


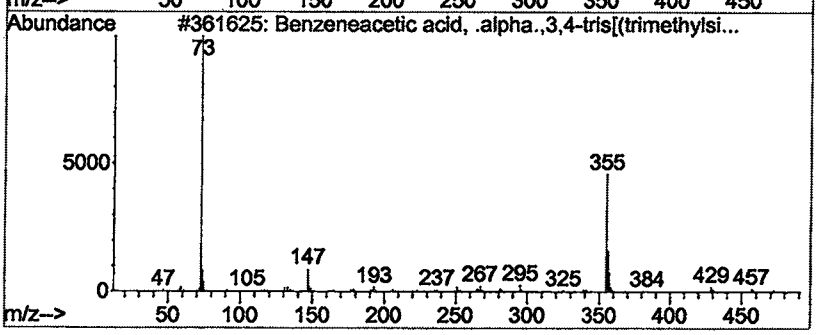


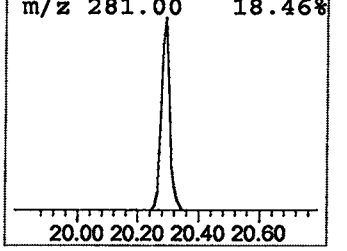












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

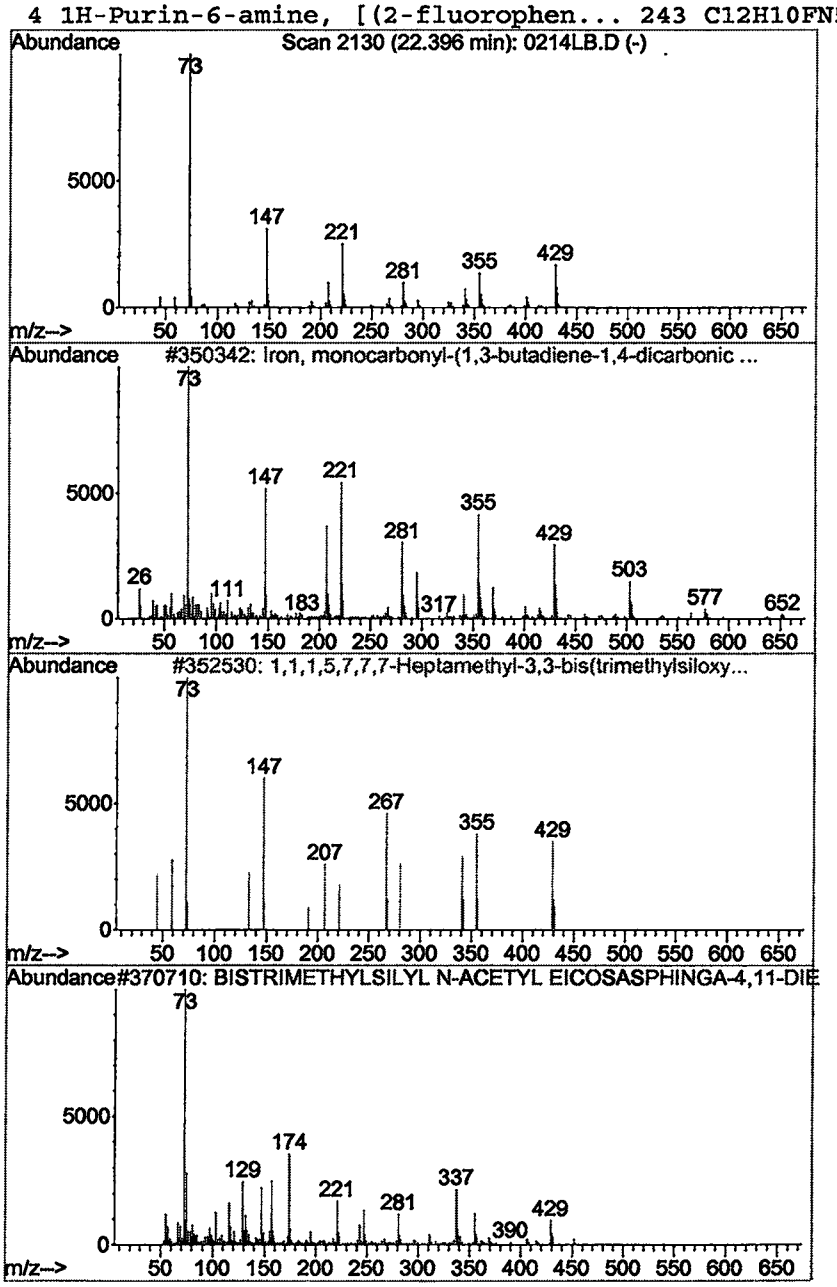
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

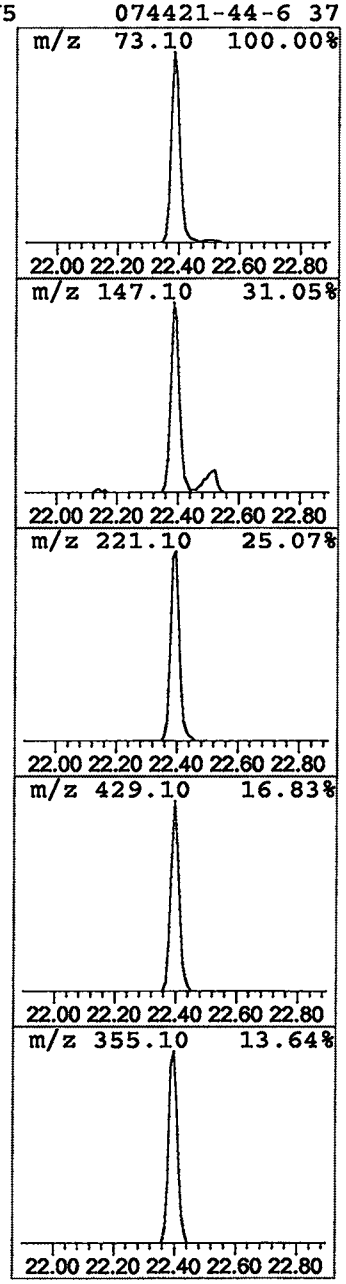
****

Peak Number 6 Iron, monocarbonyl-(1,3-but... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.40	0.81 ug/L	970014	Phenanthrene-d10	22.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Iron, monocarbonyl-(1,3-butadien...	438	C21H22FeN2O5	109007-87-6	83
2	1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	40
3	BISTRIMETHYLSILYL N-ACETYL EICOS...	511	C28H57NO3Si2	000000-00-0	40
4	1H-Purin-6-amine, [(2-fluorophen...	243	C12H10FN5	074421-44-6	37





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

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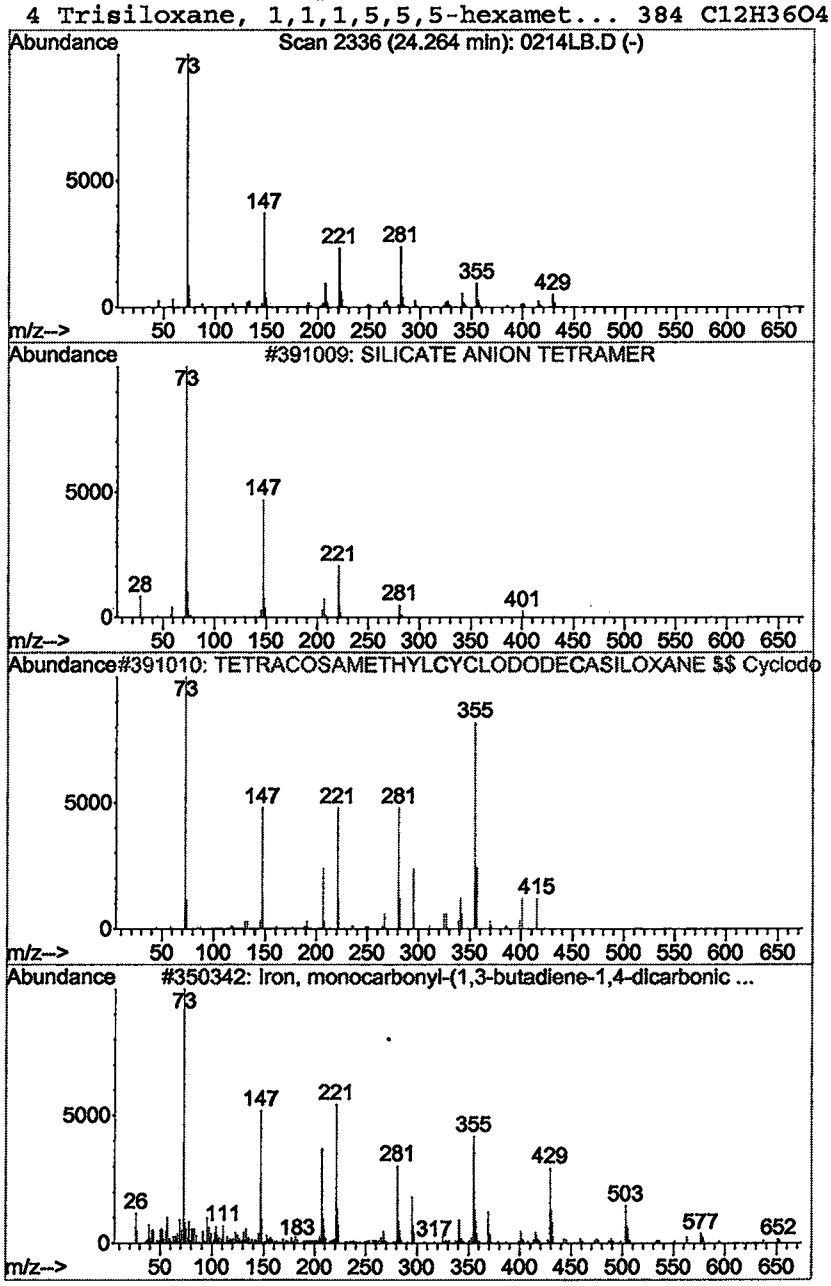
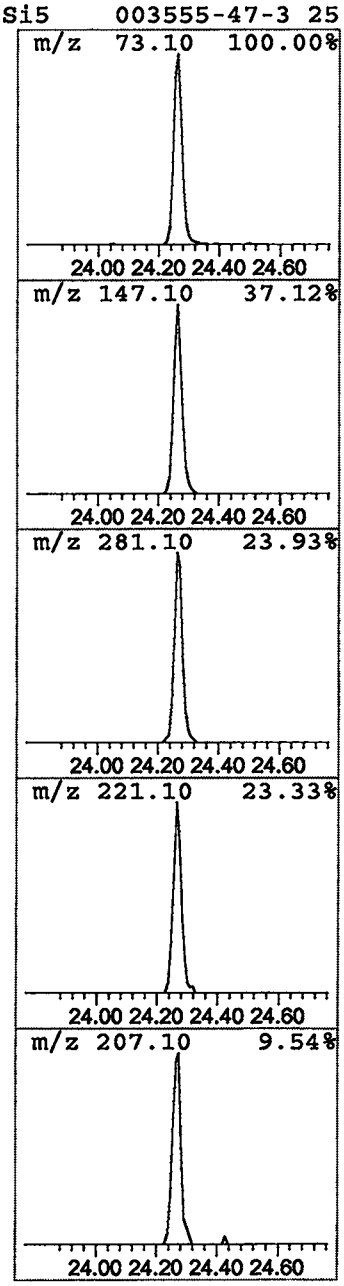
Peak Number 7 SILICATE ANION TETRAMER Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.26	0.61 ug/L	727692	Phenanthrene-d10	22.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	SILICATE ANION TETRAMER	888	C24H72O12Si12	000000-00-0	47
2	TETRACOSAMETHYLCYCLODODECASILOXA...	888	C24H72O12Si12	018919-94-3	40
3	Iron, monocarbonyl-(1,3-butadien...	438	C21H22FeN2O5	109007-87-6	35
4	Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	25





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

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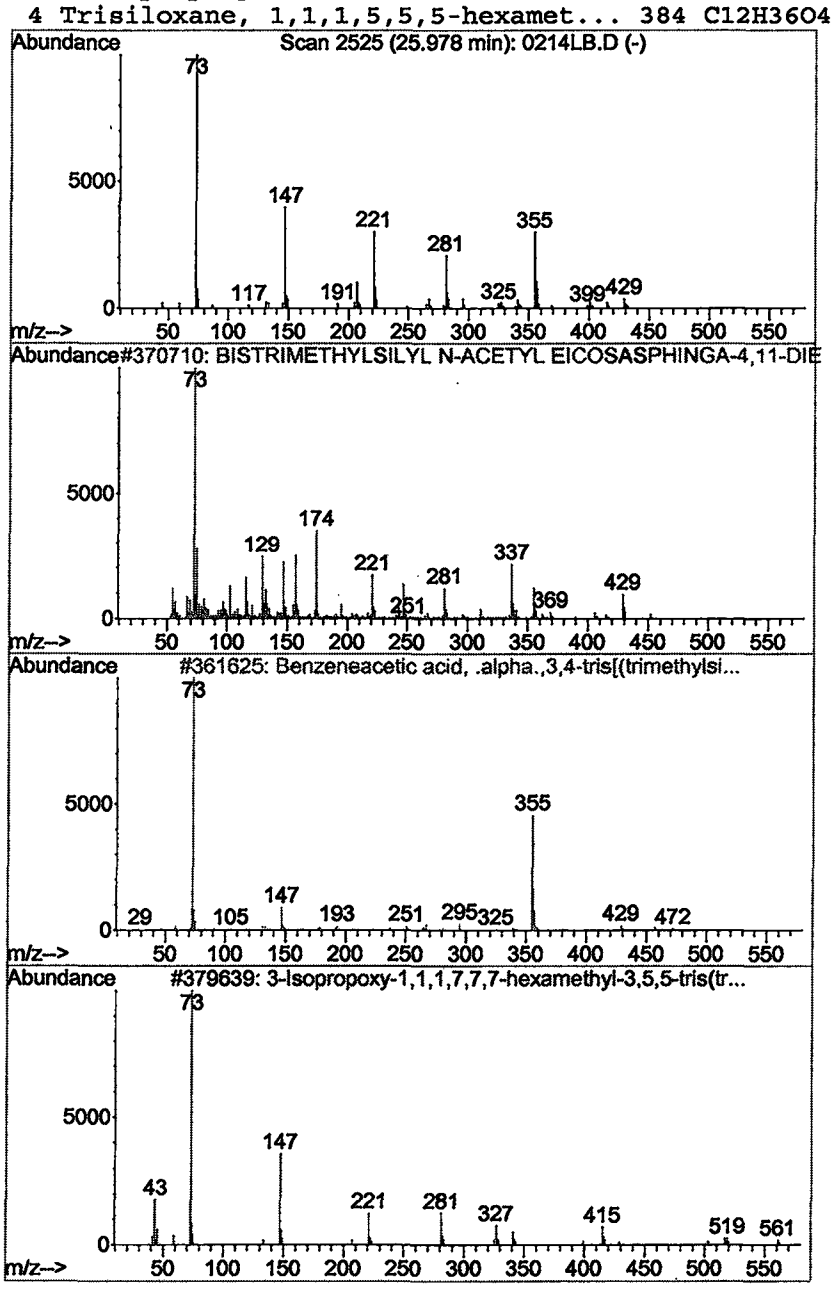
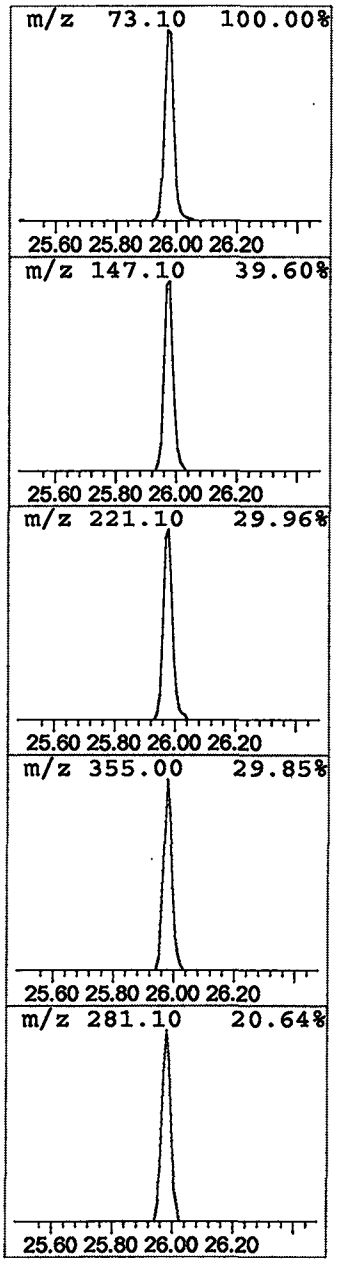
Peak Number 8 BISTRIMETHYLSILYL N-ACETYL ... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.98	0.53 ug/L	640154	Phenanthrene-d10	22.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	BISTRIMETHYLSILYL N-ACETYL EICOS...	511	C28H57NO3Si2	000000-00-0	59
2	Benzeneacetic acid, .alpha.,3,4-...	472	C20H40O5Si4	037148-65-5	40
3	3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	37
4	Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	37





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

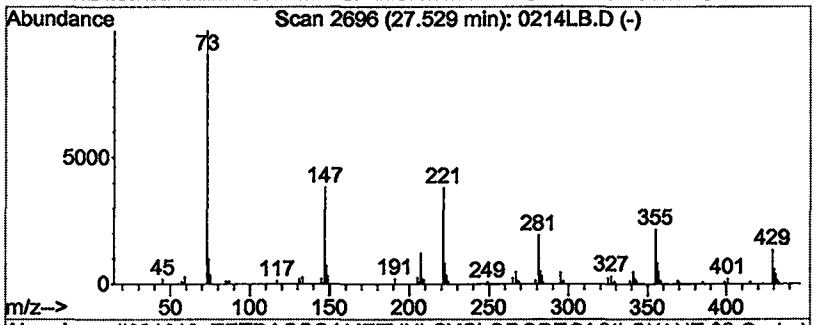
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

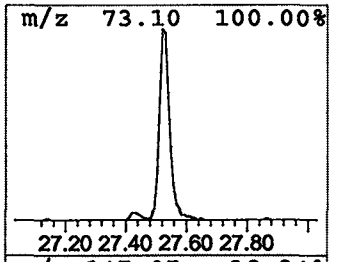
****

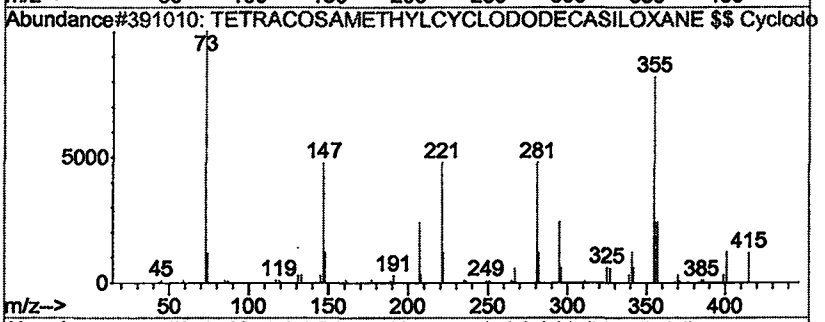
Peak Number 9 TETRACOSAMETHYLCYCLODODECAS... Concentration Rank 8

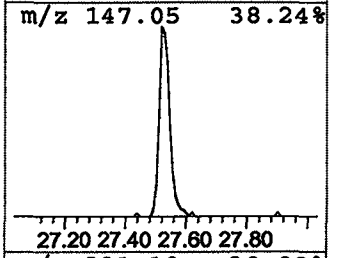
R.T.	EstConc	Area	Relative to ISTD	R.T.
27.53	0.61 ug/L	683674	Chrysene-d12	30.34

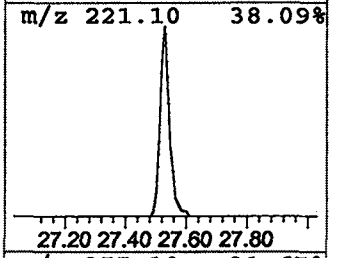
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	TETRACOSAMETHYLCYCLODODECASILOXA...	888	C24H72O12Si12	018919-94-3	76
2	1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	38
3	SILICATE ANION TETRAMER	888	C24H72O12Si12	000000-00-0	35
4	BISTRIMETHYLSILYL N-ACETYL EICOS...	511	C28H57NO3Si2	000000-00-0	27

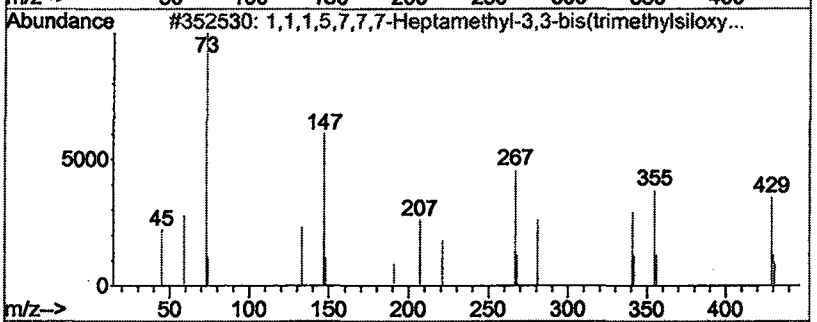


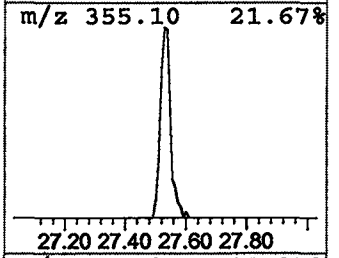


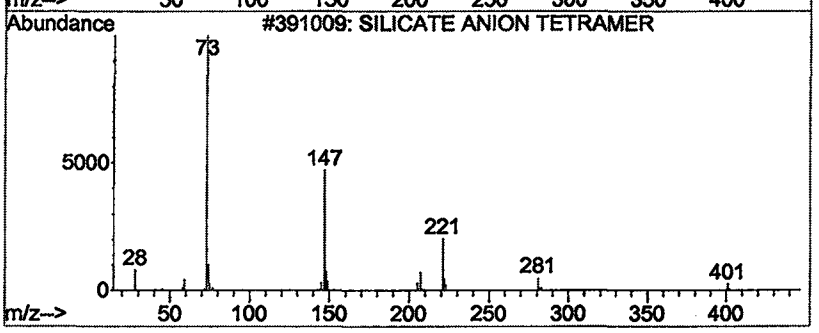


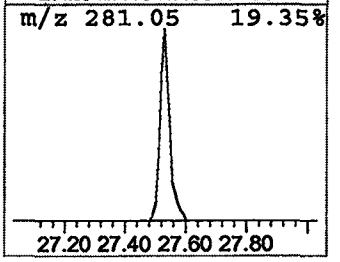












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\031902\0214LB.D Vial: 33
Acq On : 21 Mar 2002 5:58 am Operator: McLean
Sample : 2/14/02 RE-INJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

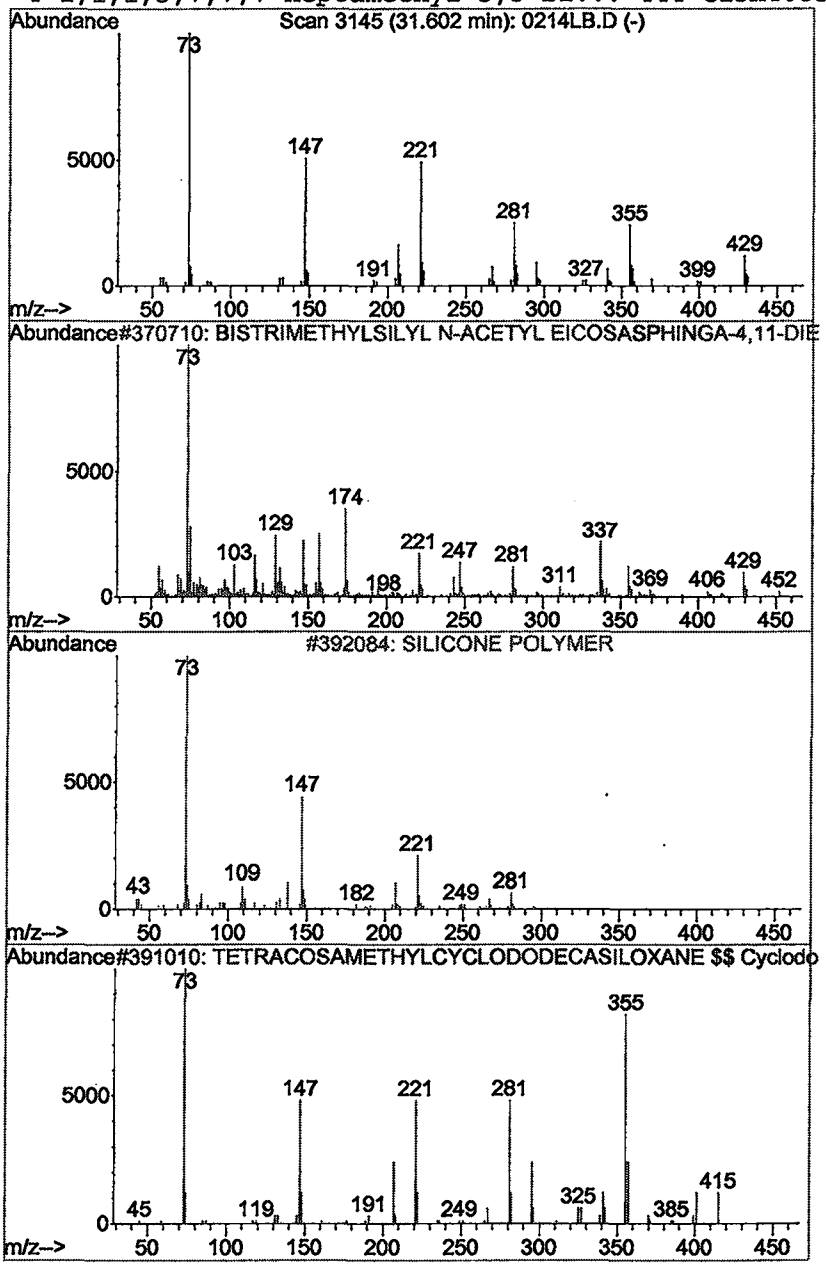
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

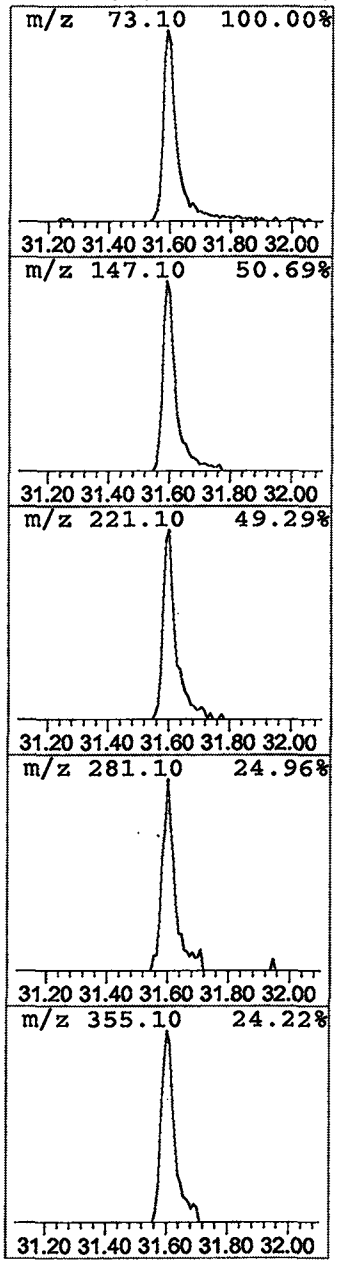
****

Peak Number 10 BISTRIMETHYLSILYL N-ACETYL ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.60	0.44 ug/L	498705	Chrysene-d12	30.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	BISTRIMETHYLSILYL N-ACETYL EICOS...	511	C28H57NO3Si2	000000-00-0	43
2	SILICONE POLYMER	458	C14H42O5Si6	000000-00-0	43
3	TETRACOSAMETHYLCYCLODODECASILOXA...	888	C24H72O12Si12	018919-94-3	38
4	1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	28





Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 21 Mar 2002 5:58 am
Data File: C:\MSDCHEM\1\DATA\031902\0214LB.D
Name: 2/14/02 RE-INJ
Misc:
Method: C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title: Quantitation For Method 508gcscan
Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	Internal Standard #	Internal Standard RT	Internal Standard Resp	Internal Standard Conc
Silane, tetramethyl-	5.56	2.2	ug/L	1833320	1	9.55	16543100	20.0
Cyclopentasiloxan...	12.06	0.7	ug/L	748627	2	13.04	22057700	20.0
Cyclohexasiloxane...	15.11	1.7	ug/L	1849620	2	13.04	22057700	20.0
Trisiloxane, 1,1,...	17.85	1.3	ug/L	1524380	3	18.15	23282000	20.0
Phenethylamine, N...	20.29	1.0	ug/L	1202490	3	18.15	23282000	20.0
Iron, monocarbony...	22.40	0.8	ug/L	970014	4	22.51	24010800	20.0
SILICATE ANION TE...	24.26	0.6	ug/L	727692	4	22.51	24010800	20.0
BISTRIMETHYLSILYL...	25.98	0.5	ug/L	640154	4	22.51	24010800	20.0
TETRACOSAMETHYLCY...	27.53	0.6	ug/L	683674	5	30.34	22566600	20.0
BISTRIMETHYLSILYL...	31.60	0.4	ug/L	498705	5	30.34	22566600	20.0