User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/22/2002 Time: 12:29:03

Sample: AE02381
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/22/02 12:28 Ended: 3/22/02 12:28 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		89		5.0

Comments for Sample AE02381 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 03-22-2002 Time: 12:29:25

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\030402\2381.D Vial: 16
Acq On : 4 Mar 2002 10:32 pm Operator: McLean
Sample : 1/31 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Mar 18 08:08:29 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	9.58	150	1474392	20.00	ug/L	0.03
10) Naphthalene-d8	13.05	136	3827699	20.00	ug/L	0.03
17) Acenaphthene-d10	18.16	164	2173070	20.00	ug/L	0.02
29) Phenanthrene-d10	22.52	188	3317728	20.00	ug/L	0.03
38) Chrysene-d12	30.34	240	2927793	20.00	ug/L	0.02
44) Perylene-d12	34.21	264	1803517	20.00	ug/L	0.03
System Monitoring Compounds						
37) 2,4-ddd surr	27.46	235	176001	4.45	ug/L	0.03
Spiked Amount 5.000			Recovery = 89.00%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	0.00	74	0	N.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.		
34) Anthracene	0.00	178	0	N.D.		
35) Di-n-butylphthalate	0.00	149	0	N.D.	d	
36) Fluoranthene	0.00	202	0	N.D.		
39) Pyrene	0.00	202	0	N.D.		
40) Butylbenzylphthalate	0.00	149	0	N.D.		
41) Benzo (a) antracene	0.00	228	0	N.D.	d	
42) Bis (2-ethylhexyl) phthala	0.00	149	0	N.D.	d	
43) Chrysene	0.00	228	0	N.D.	d	
45) Di-n-Octylphthalate	0.00	149	0	N.D.		
46) Benzo (b) fluoranthene	0.00	252	0	N.D.		

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

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Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Mar 18 08:08:29 2002 Quant Results File: 5080302.RES

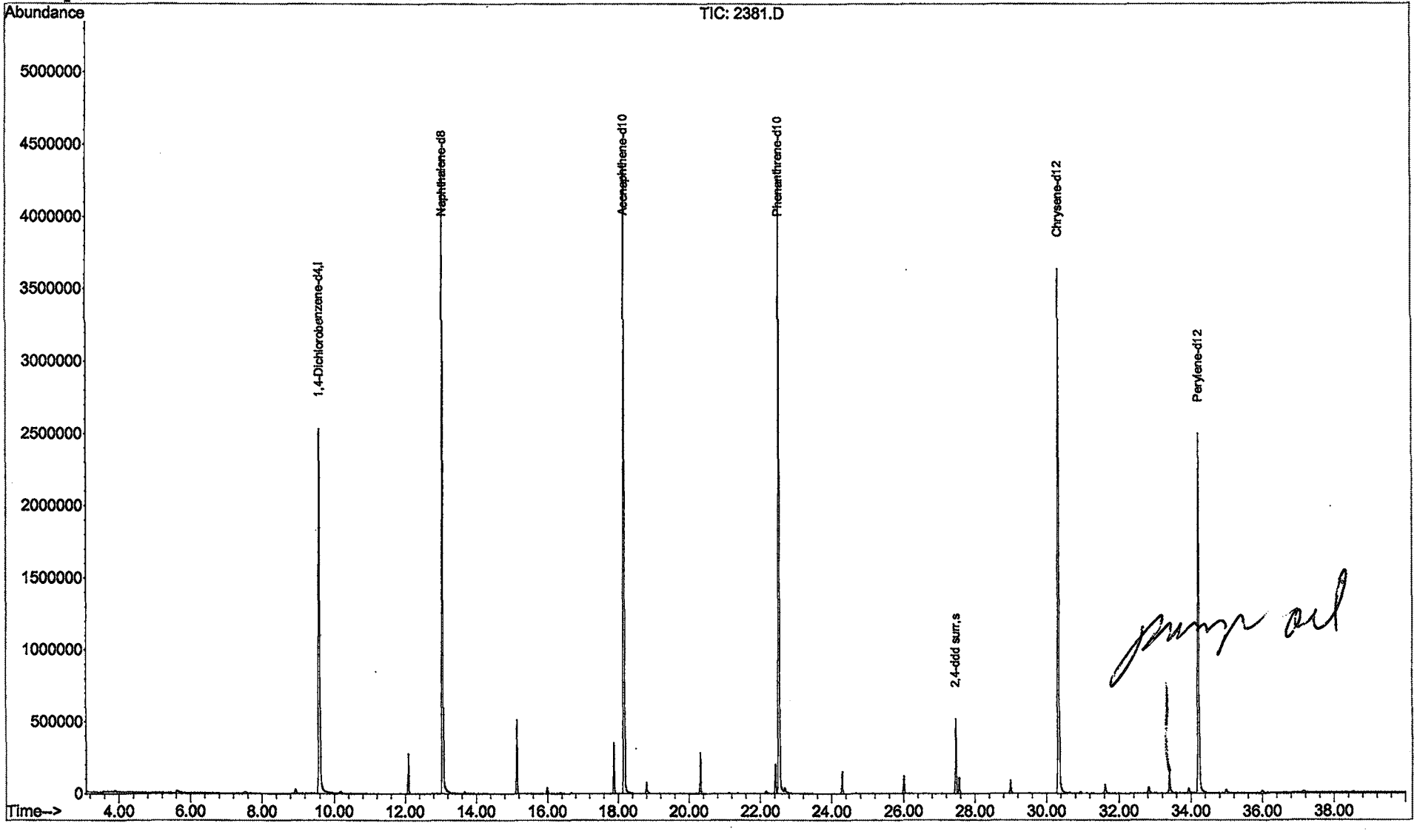
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Thu Mar 14 13:02:27 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

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
2381.D 5080302.M Mon Mar 18 08:10:18 2002 Page 2

Data File : C:\MSDCHEM\1\DATA\030402\2381.D Vial: 16
Acq On : 4 Mar 2002 10:32 pm Operator: McLean
Sample : 1/31 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Mar 18 8:09 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



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\030402\2381.D Vial: 16
Acq On : 4 Mar 2002 10:32 pm Operator: McLean
Sample : 1/31 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

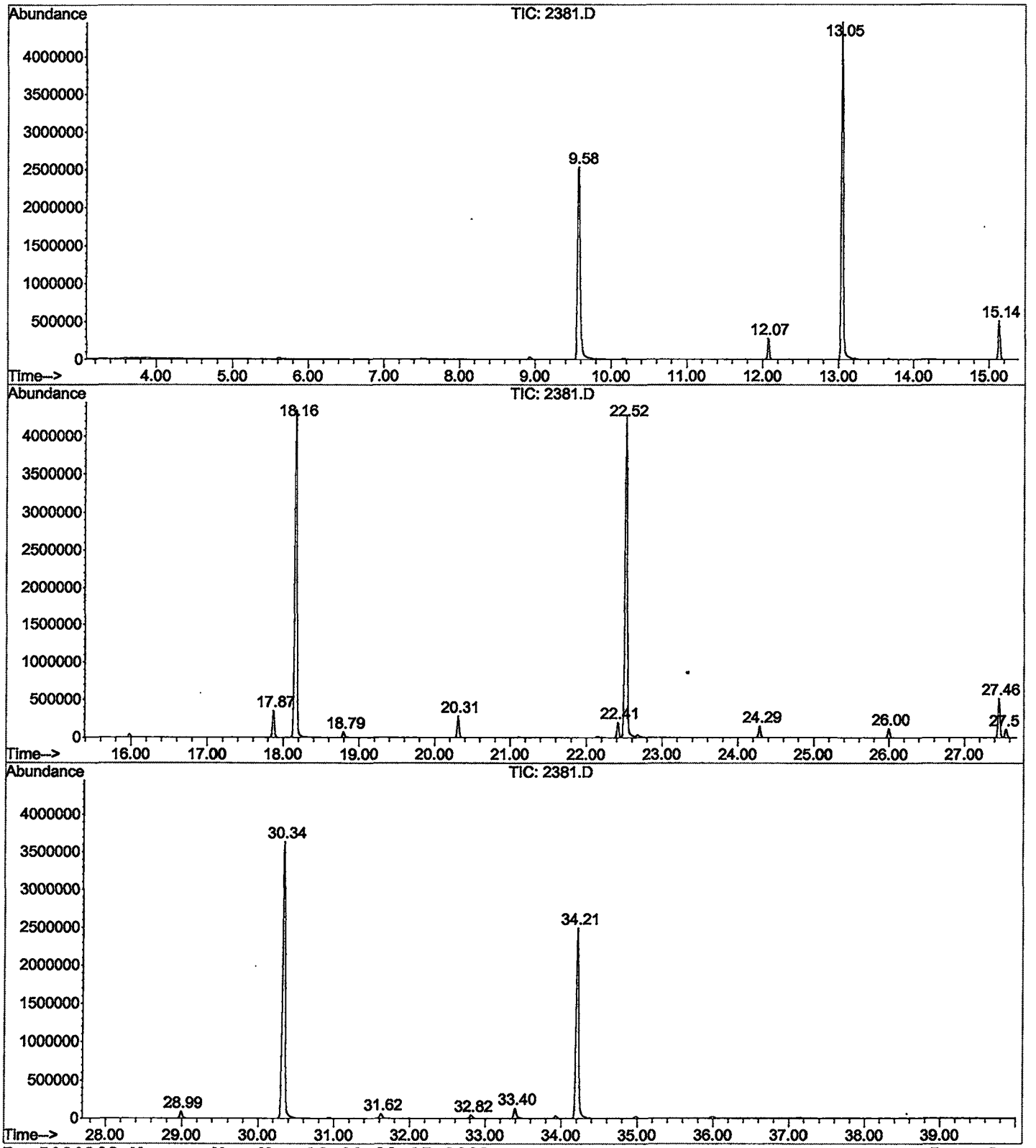
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.579	711	717	734	rBV	2532243	6327561	69.28%	11.977%
2	12.074	988	992	997	rVB	276581	464954	5.09%	0.880%
3	13.053	1094	1100	1115	rBV	4452809	8314607	91.04%	15.739%
4	15.140	1325	1330	1337	rVB2	511123	822933	9.01%	1.558%
5	17.870	1626	1631	1636	rBB	351772	606414	6.64%	1.148%
6	18.160	1657	1663	1676	rBV	4352637	9111458	99.77%	17.247%
7	18.795	1730	1733	1740	rBV	74573	135364	1.48%	0.256%
8	20.310	1895	1900	1905	rBV	282508	463608	5.08%	0.878%
9	22.414	2127	2132	2137	rBB2	203753	361098	3.95%	0.684%
10	22.523	2137	2144	2158	rBV2	4284060	9132777	100.00%	17.287%
11	24.291	2334	2339	2344	rBV	151257	259977	2.85%	0.492%
12	25.997	2523	2527	2534	rBB	119367	221262	2.42%	0.419%
13	27.457	2684	2688	2694	rBV	517651	927008	10.15%	1.755%
14	27.557	2694	2699	2706	rVB	108016	220347	2.41%	0.417%
15	28.990	2852	2857	2865	rBV	89921	177622	1.94%	0.336%
16	30.341	2998	3006	3022	rBV2	3637024	8719334	95.47%	16.505%
17	31.620	3142	3147	3155	rBV	60632	135462	1.48%	0.256%
18	32.817	3273	3279	3289	rVB	43492	102896	1.13%	0.195%
19	33.398	3338	3343	3353	rBV	134535	319809	3.50%	0.605%
20	34.214	3425	3433	3454	rBV	2496231	6004925	65.75%	11.367%

Sum of corrected areas: 52829416

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\030402\2381.D
Operator : McLean
Acquired : 4 Mar 2002 10:32 pm using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 1/31
Misc Info :
Vial Number: 16
Quant File :5080302.RES (RTE Integrator)



Library Search Compound Report

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Acq On : 4 Mar 2002 10:32 pm Operator: McLean
Sample : 1/31 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

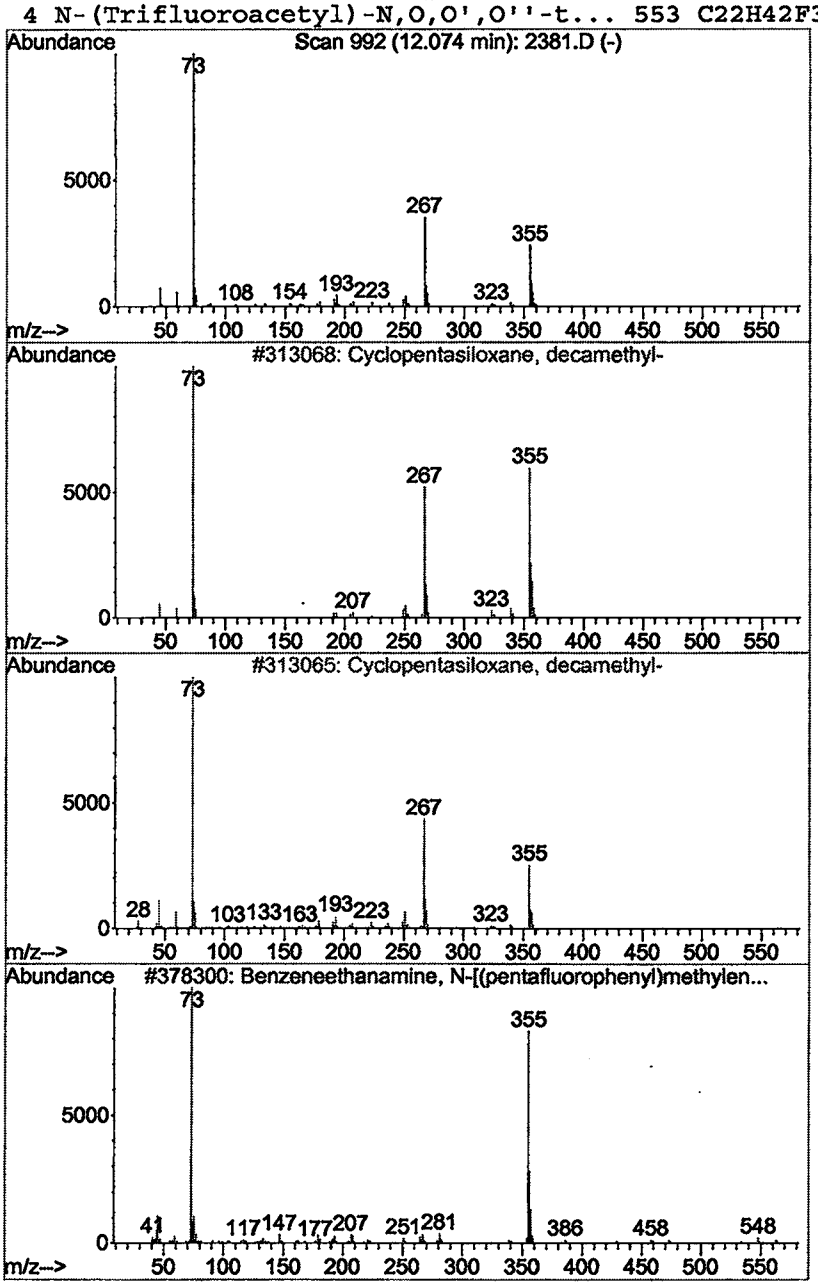
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

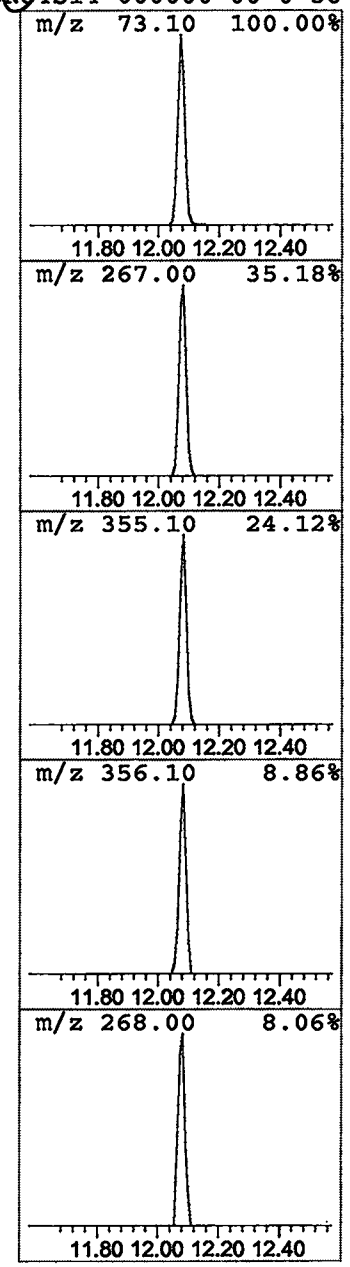
****

Peak Number 1 Cyclopentasiloxane, decamet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.07	1.12 ug/L	464954	Naphthalene-d8	13.05

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	76
3	Benzeneethanamine, N-[(pentafluo...	563	C24H34F5NO3Si3	055429-13-5	43
4	N-(Trifluoroacetyl)-N,O,O',O''-t...	553	C22H42F3NO4Si4	000000-00-0	38





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Sample : 1/31 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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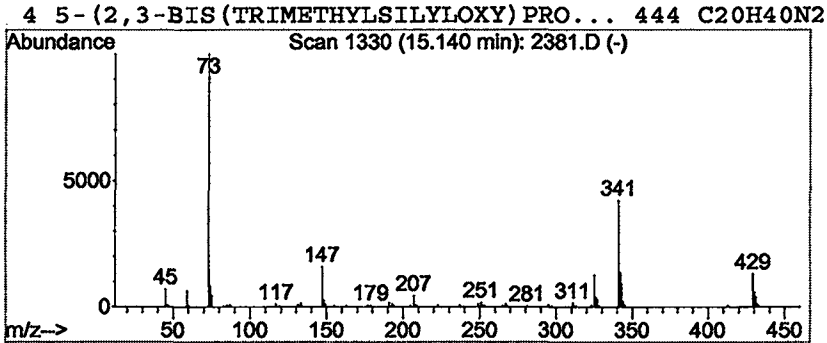
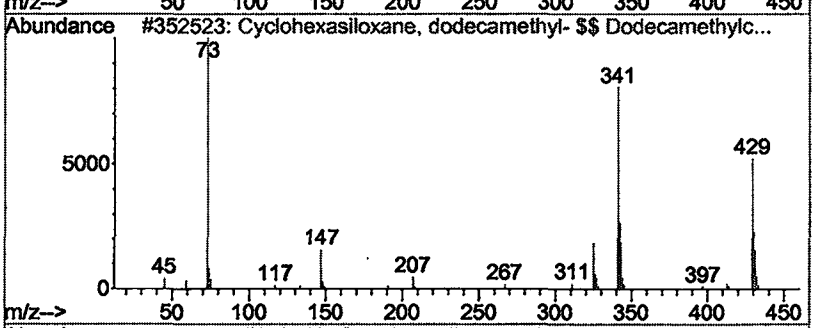
****

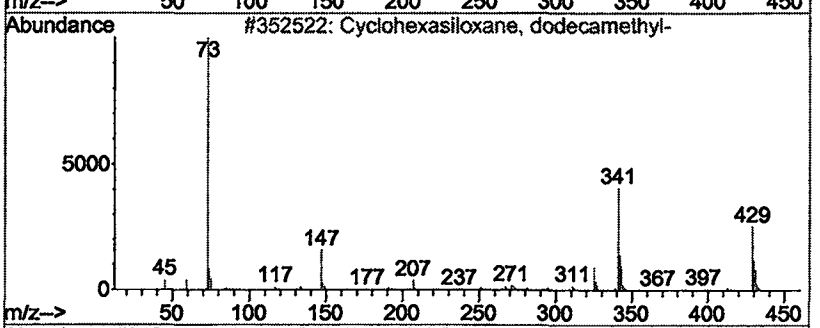
Peak Number 2 Cyclohexasiloxane, dodecame... Concentration Rank 1

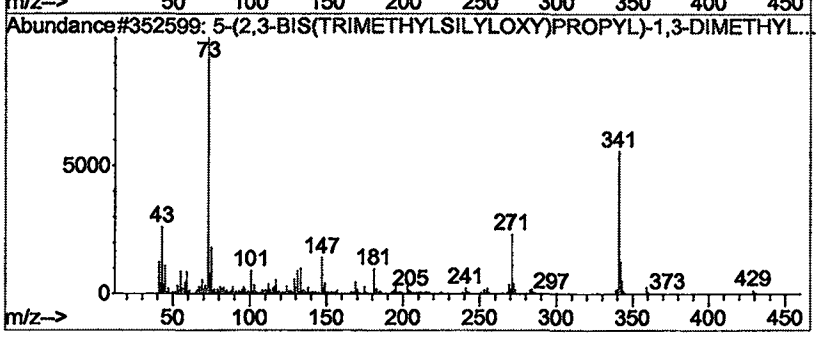
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	1.98 ug/L	822933	Naphthalene-d8	13.05

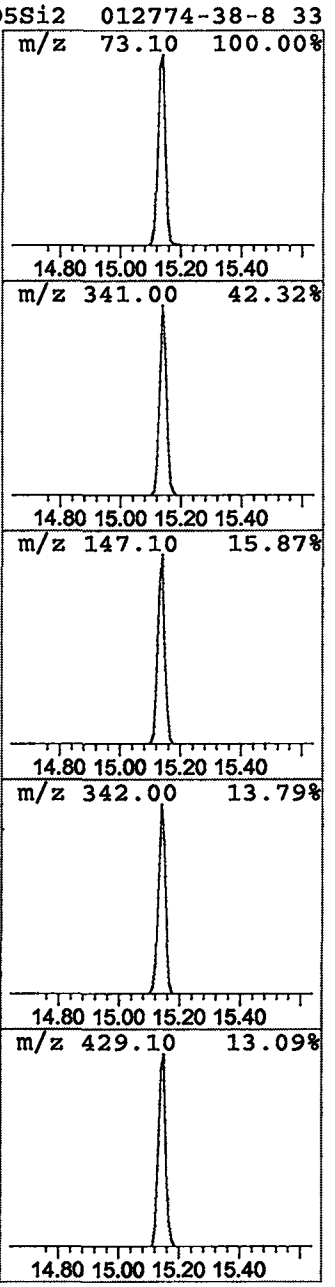
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexasiloxane, dodecamethyl-...	444	C12H36O6Si6	000540-97-6	94
2	Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	91
3	5-(2,3-BIS(TRIMETHYLSILYLOXY)PRO...	444	C20H40N2O5Si2	012774-38-8	36
4	5-(2,3-BIS(TRIMETHYLSILYLOXY)PRO...	444	C20H40N2O5Si2	012774-38-8	33











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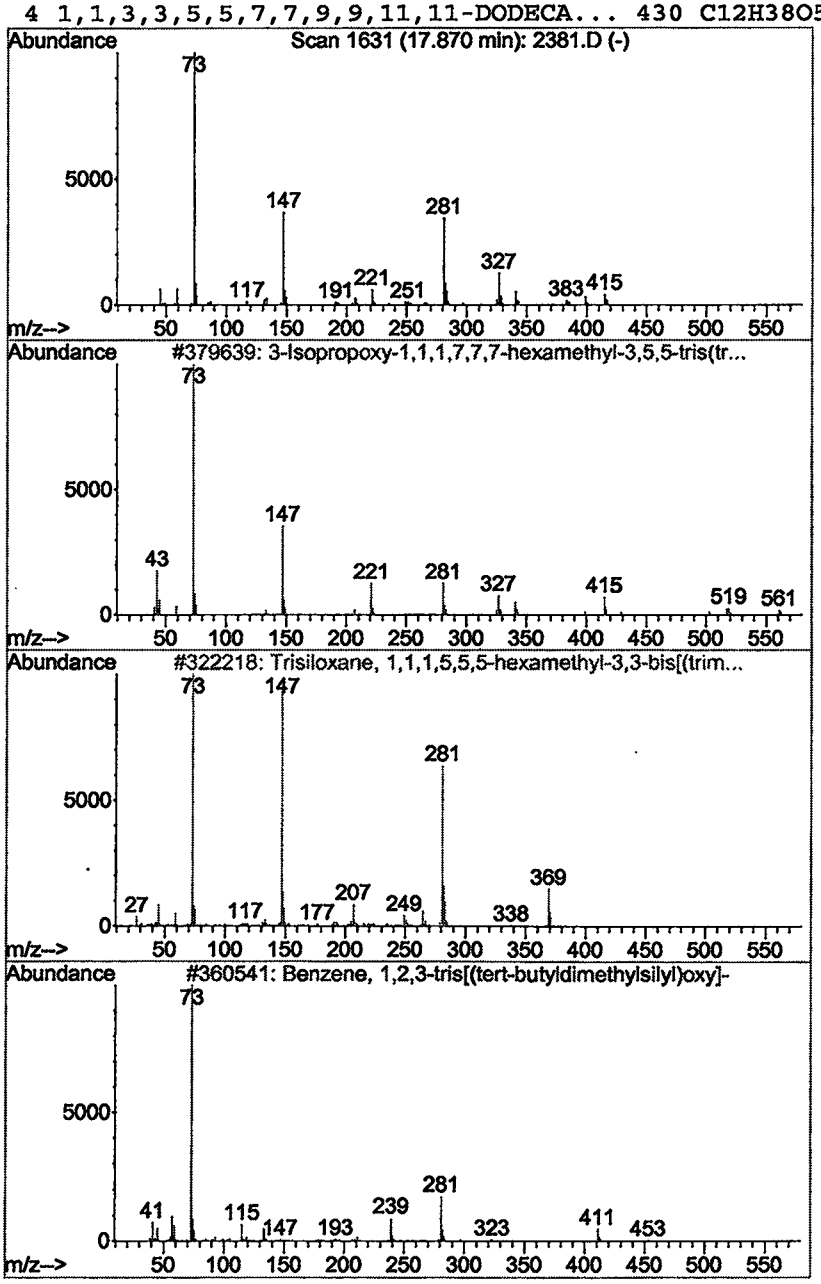
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

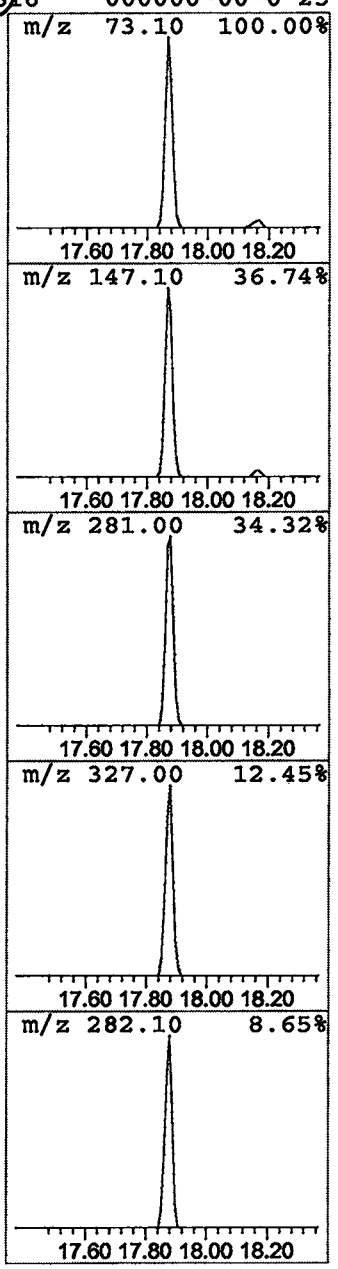
****

Peak Number 3 3-Isopropoxy-1,1,1,7,7,7-he... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.87	1.33 ug/L	606414	Acenaphthene-d10	18.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	45
2	Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	42
3	Benzene, 1,2,3-tris[(tert-butyld...	468	C24H48O3Si3	000000-00-0	35
4	1,1,3,3,5,5,7,7,9,9,11,11-DODECA...	430	C12H38O5Si6	000000-00-0	25





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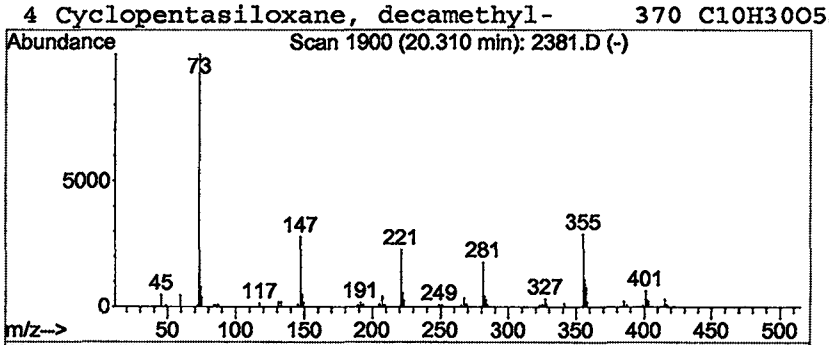
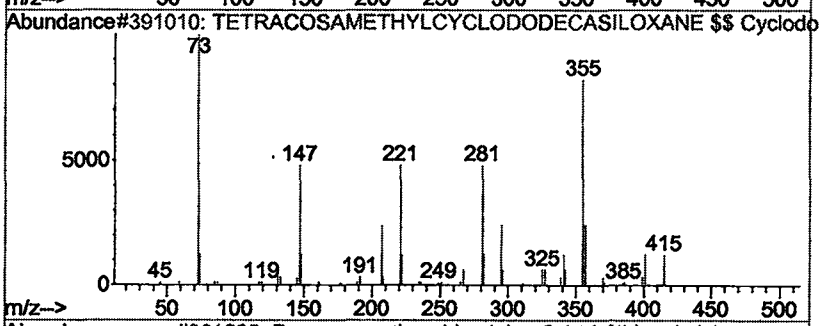
****

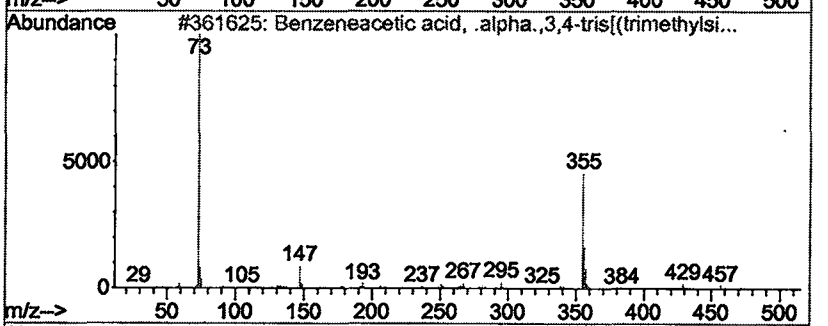
Peak Number 4 TETRACOSAMETHYLCYCLODODECAS... Concentration Rank 5

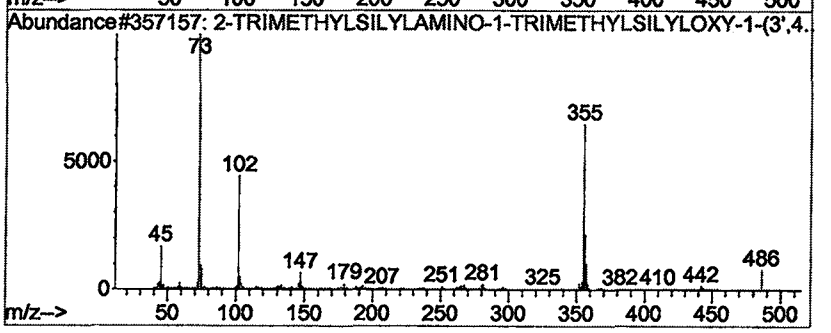
R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	1.02 ug/L	463608	Acenaphthene-d10	18.16

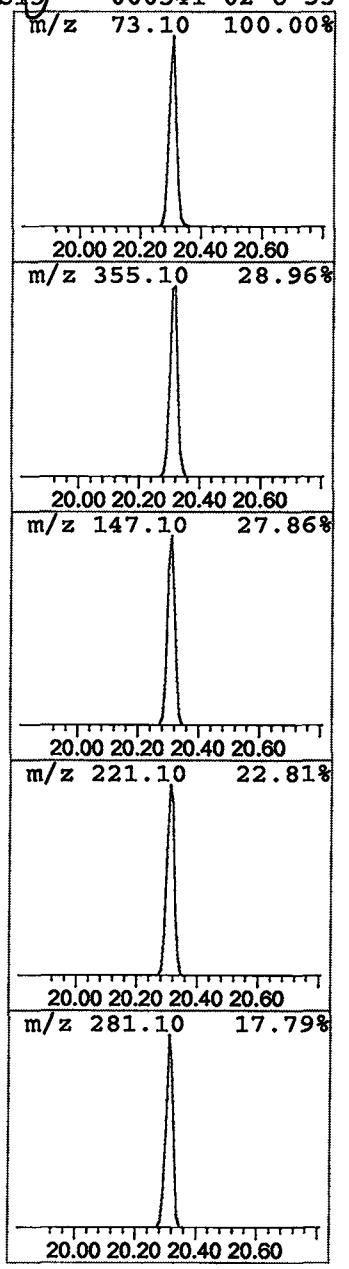
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	TETRACOSAMETHYLCYCLODODECASILOXA...	888	C24H72O12Si12	018919-94-3	56
2	Benzeneacetic acid, .alpha.,3,4-...	472	C20H40O5Si4	037148-65-5	40
3	2-TRIMETHYLSILYLAMINO-1-TRIMETHY...	457	C20H43NO3Si4	068595-65-3	35
4	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	35











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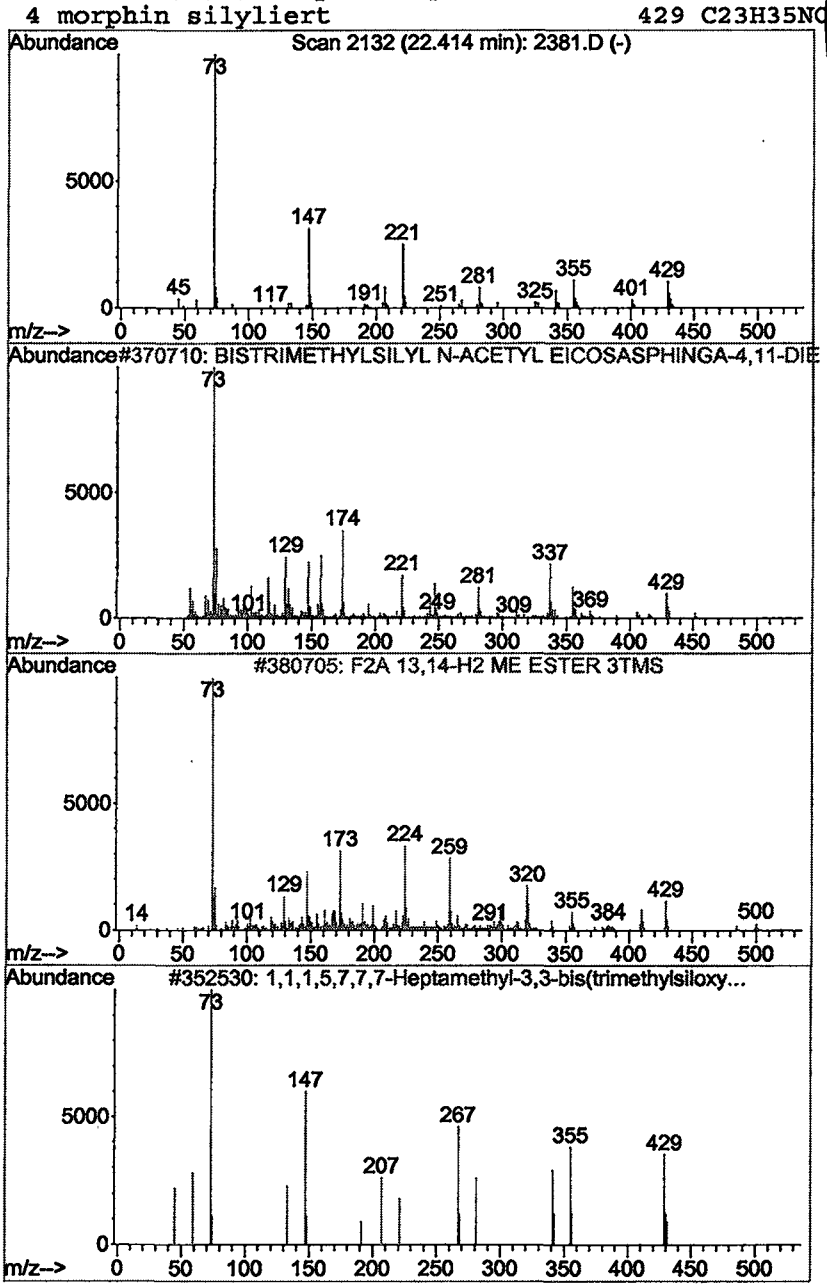
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

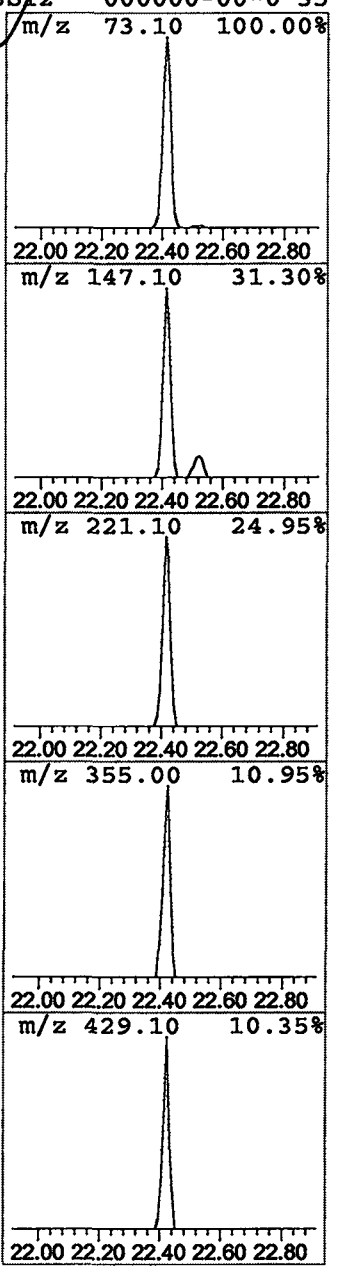
****

Peak Number 5 BISTRIMETHYLSILYL N-ACETYL ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.41	0.79 ug/L	361098	Phenanthrene-d10	22.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	BISTRIMETHYLSILYL N-ACETYL EICOS...	511	C28H57NO3Si2	000000-00-0	47
2	F2A 13,14-H2 ME ESTER 3TMS	586	C30H62O5Si3	000000-00-0	42
3	1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	40
4	morphin silyliert	429	C23H35NO3Si2	000000-00-0	35





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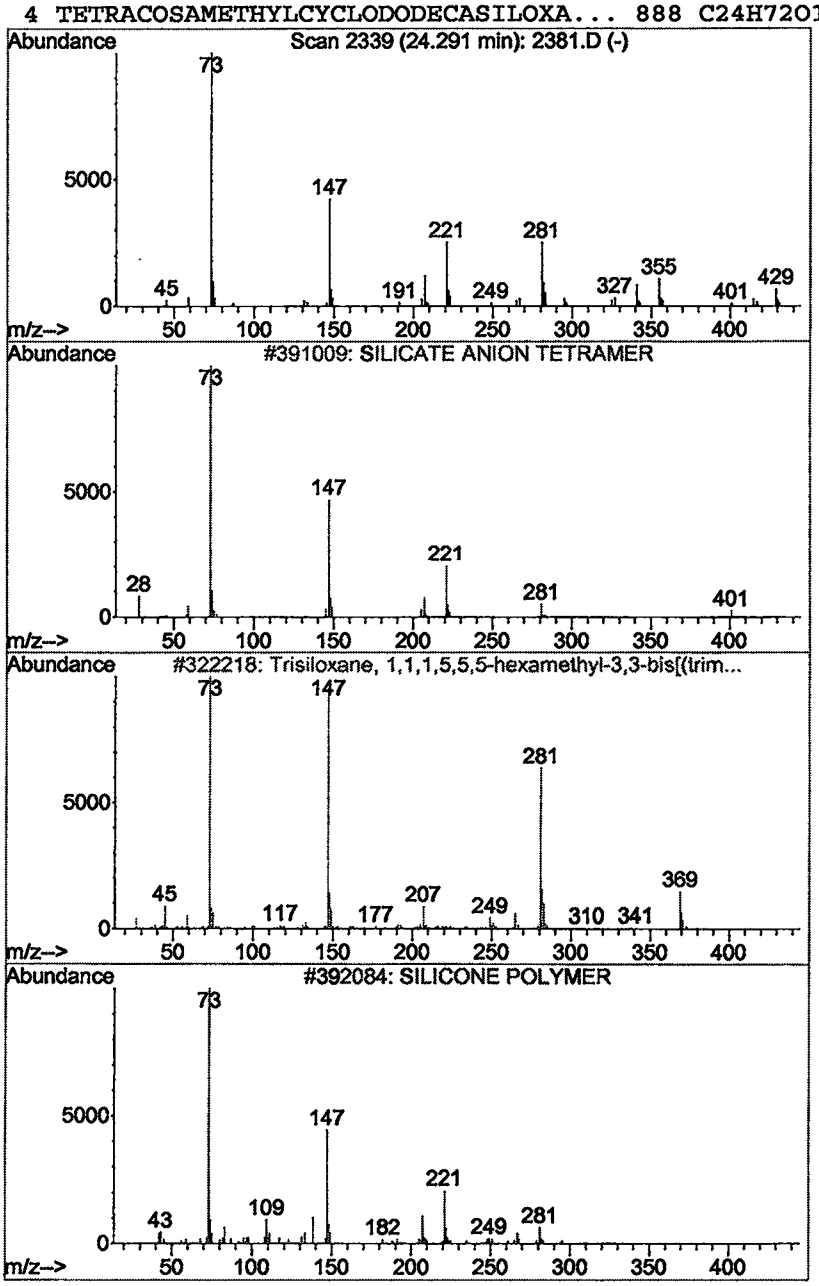
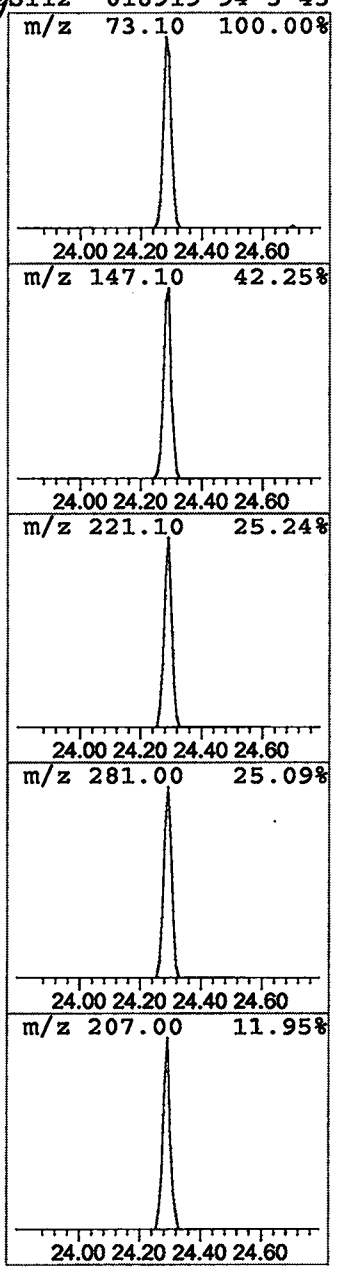
****

Peak Number 6 SILICATE ANION TETRAMER Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.29	0.57 ug/L	259977	Phenanthrene-d10	22.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	SILICATE ANION TETRAMER	888	C24H72O12Si12	000000-00-0	50
2	Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	50
3	SILICONE POLYMER	458	C14H42O5Si6	000000-00-0	43
4	TETRACOSAMETHYLCYCLODODECASILOXA...	888	C24H72O12Si12	018919-94-3	43





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Data File : C:\MSDCHEM\1\DATA\030402\2381.D Vial: 16
Acq On : 4 Mar 2002 10:32 pm Operator: McLean
Sample : 1/31 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

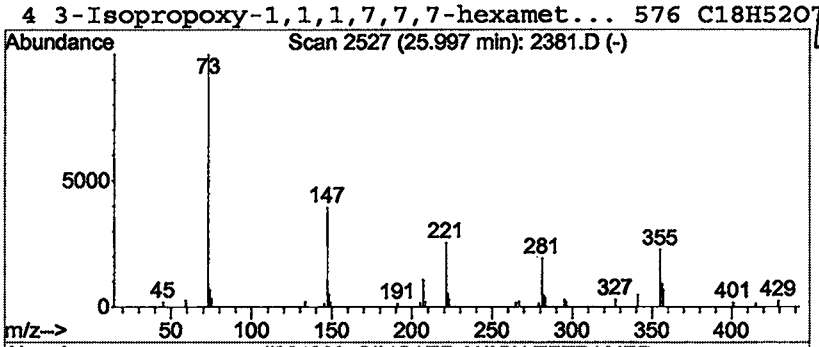
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

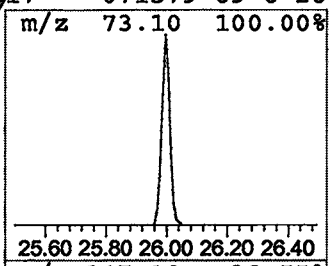
****

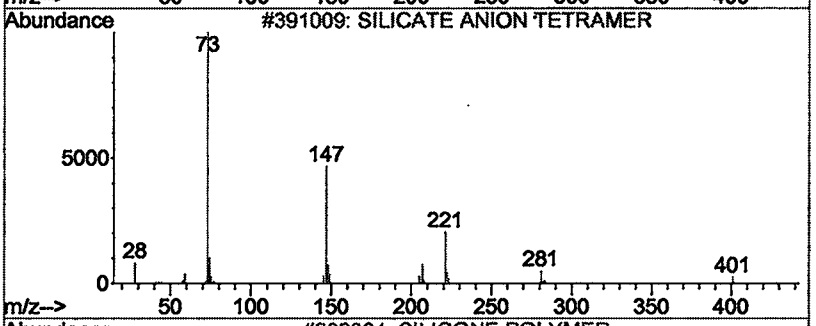
Peak Number 7 SILICATE ANION TETRAMER Concentration Rank 9

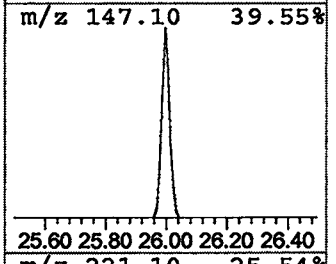
R.T.	EstConc	Area	Relative to ISTD	R.T.
26.00	0.48 ug/L	221262	Phenanthrene-d10	22.52

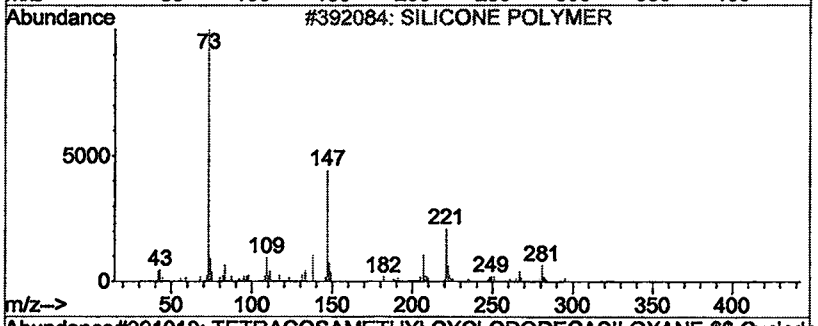
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	SILICATE ANION TETRAMER	888	C24H72O12Si12	000000-00-0	38
2	SILICONE POLYMER	458	C14H42O5Si6	000000-00-0	38
3	TETRACOSAMETHYLCYCLODODECASILOXA...	888	C24H72O12Si12	018919-94-3	32
4	3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	28

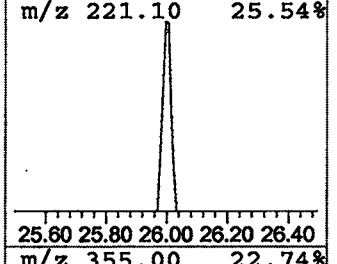


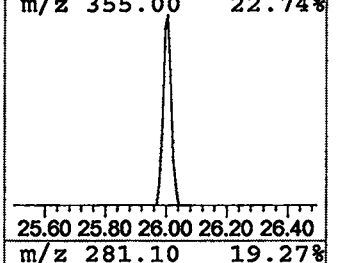


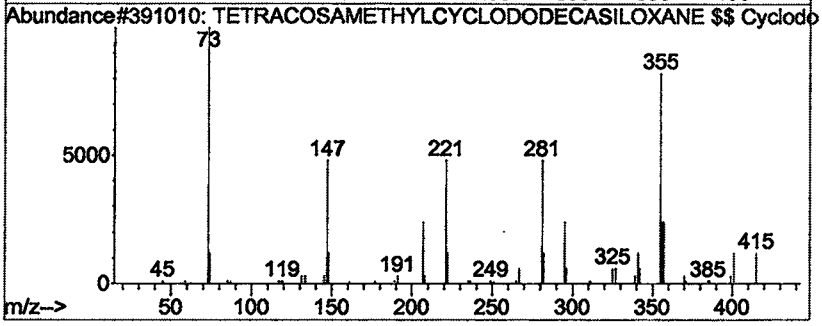


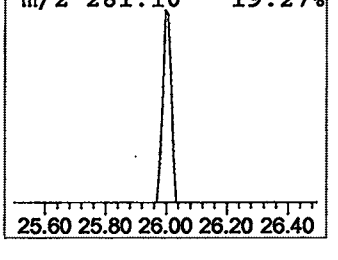












Library Search Compound Report

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Acq On : 4 Mar 2002 10:32 pm Operator: McLean
Sample : 1/31 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

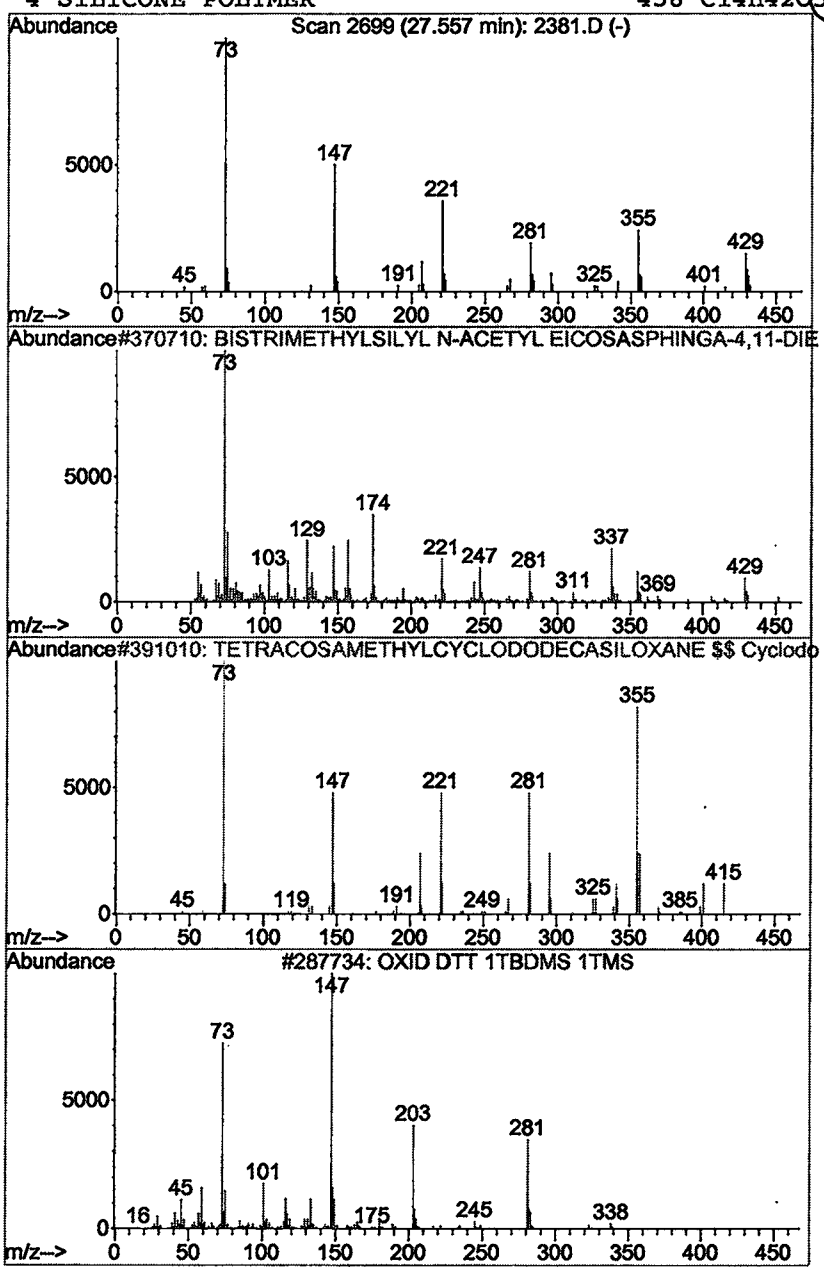
Quant Method : C:\MSDCHEM\1\5973N\5080302.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

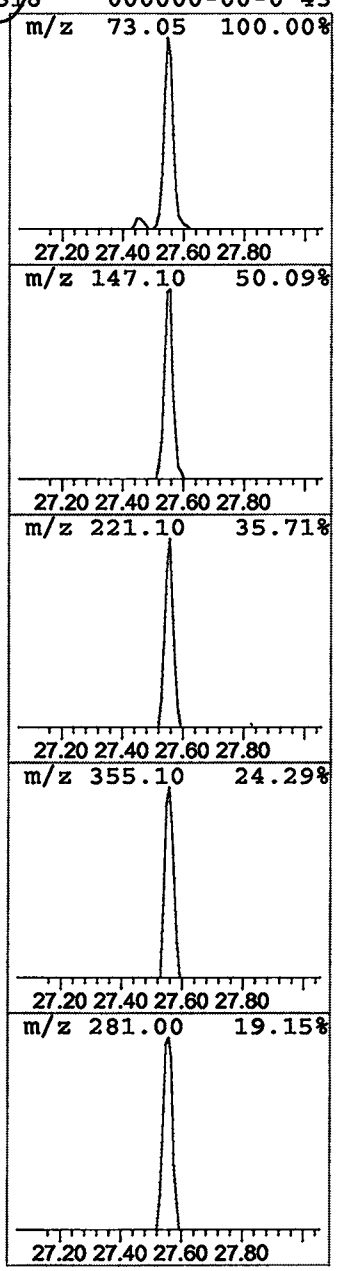
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.56	0.51 ug/L	220347	Chrysene-d12	30.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	BISTRIMETHYLSILYL N-ACETYL EICOS...	511	C28H57NO3Si2	000000-00-0	53
2	TETRACOSAMETHYLCYCLODODECASILOXA...	888	C24H72O12Si12	018919-94-3	49
3	OXID DTT 1TBDMS 1TMS	338	C13H30O2S2Si2	000000-00-0	43
4	SILICONE POLYMER	458	C14H42O5Si6	000000-00-0	43





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Sample : 1/31 Inst : Instrumen
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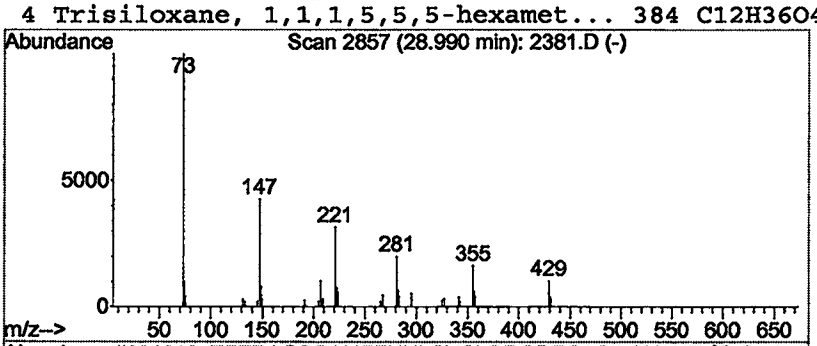
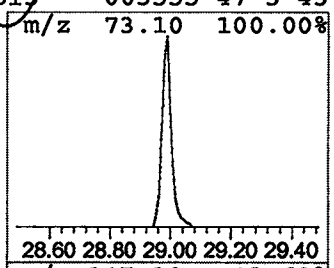
****

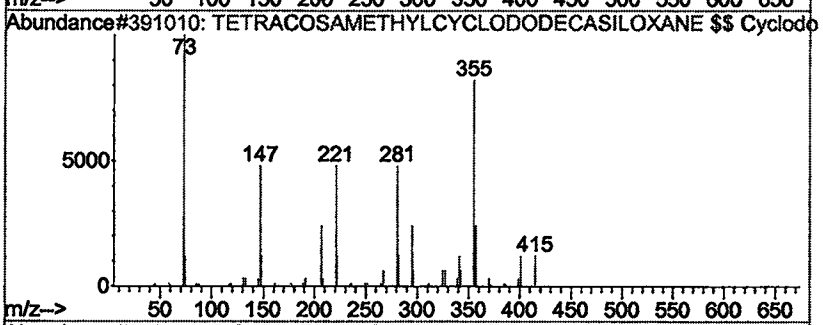
Peak Number 9 TETRACOSAMETHYLCYCLODODECAS... Concentration Rank 10

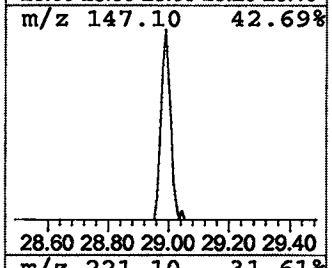
R.T.	EstConc	Area	Relative to ISTD	R.T.
28.99	0.41 ug/L	177622	Chrysene-d12	30.34

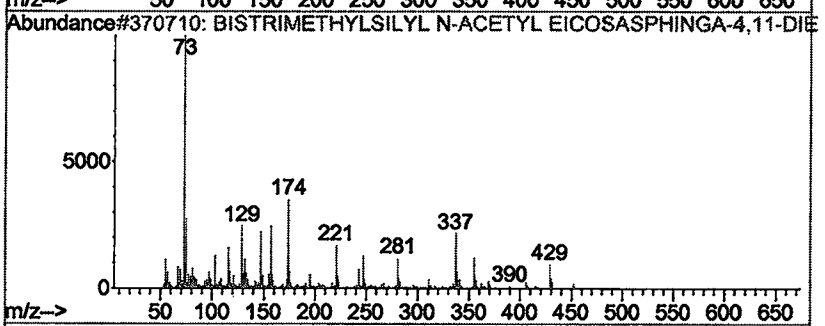
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	TETRACOSAMETHYLCYCLODODECASILOXA...	888	C24H72O12Si12	018919-94-3	83
2	BISTRIMETHYLSILYL N-ACETYL EICOS...	511	C28H57NO3Si2	000000-00-0	50
3	Iron, monocarbonyl-(1,3-butadien...	438	C21H22FeN2O5	109007-87-6	46
4	Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	43

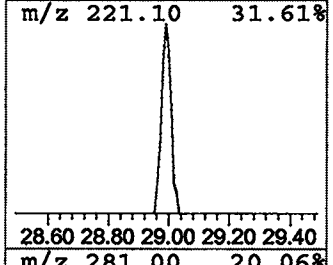


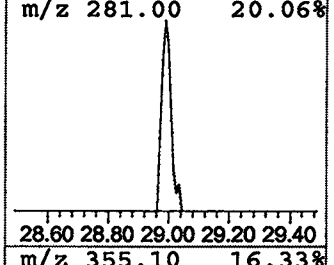


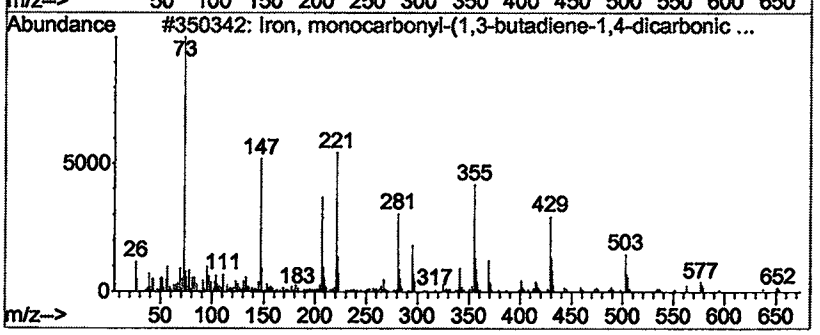


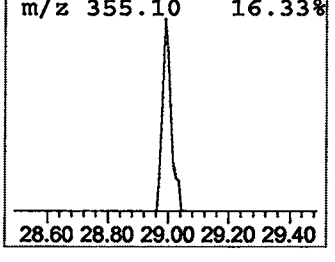












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Title : Quantitation For Method 508gcscan
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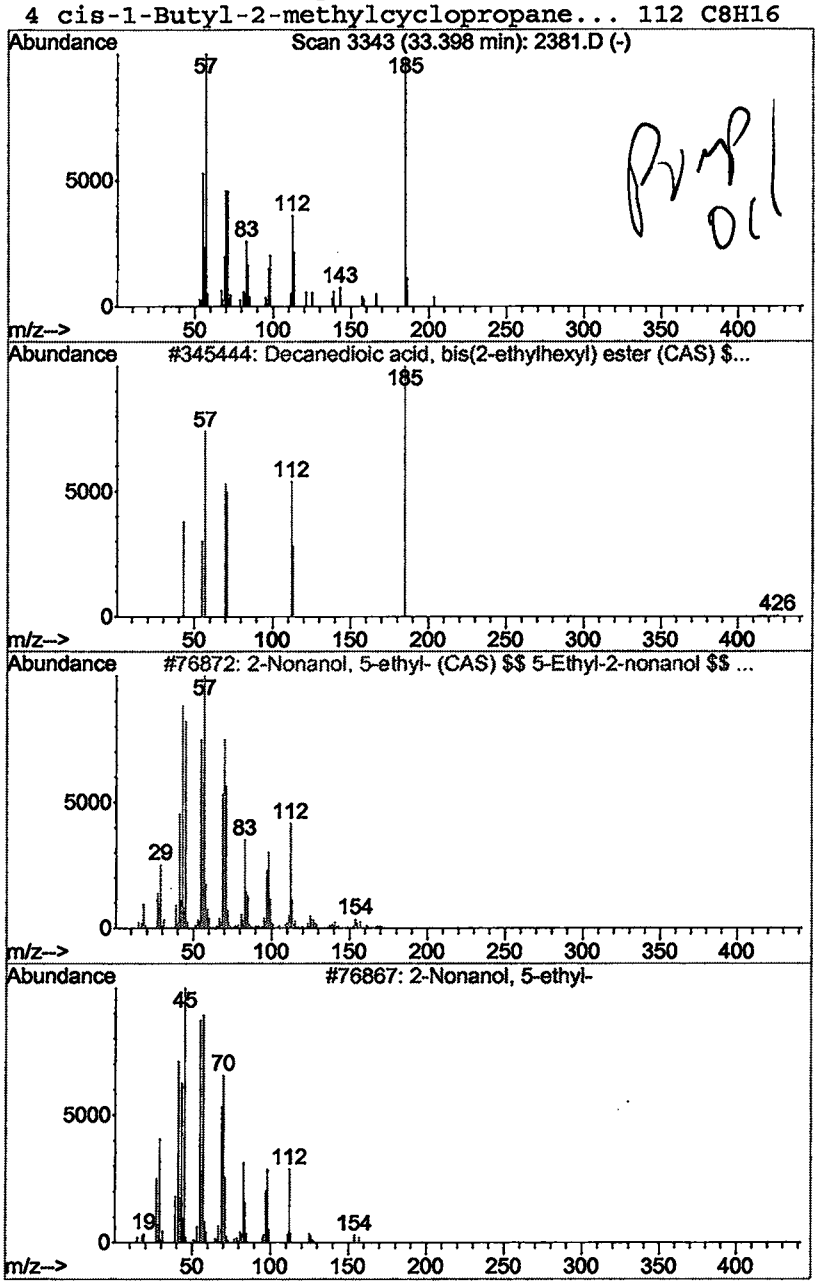
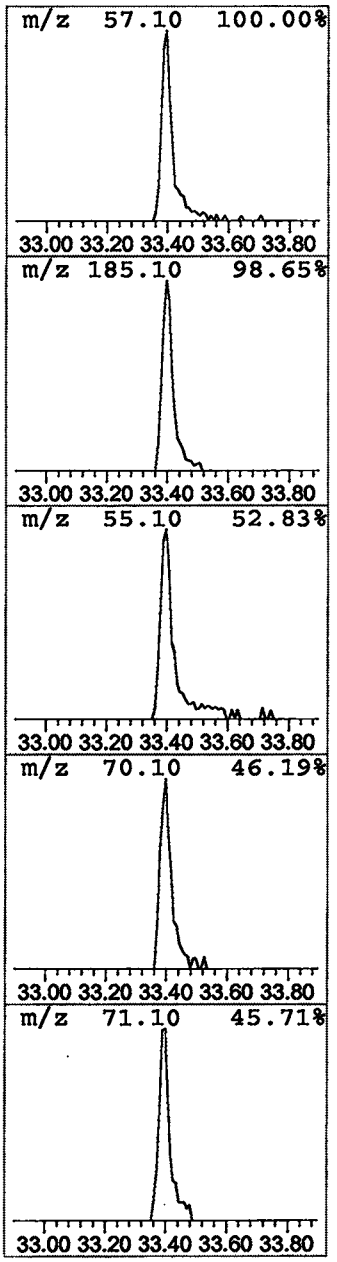
****

Peak Number 10 Decanedioic acid, bis(2-eth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
33.40	1.07 ug/L	319809	Perylene-d12	34.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decanedioic acid, bis(2-ethylhex...	426	C26H50O4	000122-62-3	59
2	2-Nonanol, 5-ethyl- (CAS) $$ 5-E...	172	C11H24O	000103-08-2	47
3	2-Nonanol, 5-ethyl-	172	C11H24O	000103-08-2	35
4	cis-1-Butyl-2-methylcyclopropane...	112	C8H16	038851-69-3	22





Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 4 Mar 2002 10:32 pm
Data File: C:\MSDCHEM\1\DATA\030402\2381.D
Name: 1/31
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
Cyclopentasiloxan...	12.07	1.1	ug/L	464954	2	13.05	8314610	20.0
Cyclohexasiloxane...	15.14	2.0	ug/L	822933	2	13.05	8314610	20.0
3-Isopropoxy-1,1,...	17.87	1.3	ug/L	606414	3	18.16	9111460	20.0
TETRACOSAMETHYLCY...	20.31	1.0	ug/L	463608	3	18.16	9111460	20.0
BISTRIMETHYLSILYL...	22.41	0.8	ug/L	361098	4	22.52	9132780	20.0
SILICATE ANION TE...	24.29	0.6	ug/L	259977	4	22.52	9132780	20.0
SILICATE ANION TE...	26.00	0.5	ug/L	221262	4	22.52	9132780	20.0
BISTRIMETHYLSILYL...	27.56	0.5	ug/L	220347	5	30.34	8719330	20.0
TETRACOSAMETHYLCY...	28.99	0.4	ug/L	177622	5	30.34	8719330	20.0
Decanedioic acid,...	33.40	1.1	ug/L	319809	6	34.21	6004930	20.0