

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 04/18/2002 Time: 12:23:10

Sample: AE06585
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 4/18/2002 12:22 Ended: 4/18/2002 12:22 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	MDL
1	Other unknown compounds		.	
2	Surr_2,4-DDD		105	5.0

Comments for Sample AE06585 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-18-2002 Time: 12:23:23

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\041502\6585.D

Vial: 9

Acq On : 15 Apr 2002 4:38 pm

Operator: Bohlk

Sample : SAMPLE 6585 3/20/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 17 16:41:14 2002

Quant Results File: 5080402BBC.RES

Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Tue Apr 16 18:27:45 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.41	152	426175	40.00	ug/L	0.00
10) Naphthalene-d8	12.88	136	1945895	40.00	ug/L	0.00
17) Acenaphthene-d10	17.99	164	1096698	40.00	ug/L	0.00
30) Phenanthrene-d10	22.34	188	1567896	40.00	ug/L	0.00
39) Chrysene-d12	30.14	240	1328134	40.00	ug/L	-0.01
45) Perylene-d12	34.01	264	830365	40.00	ug/L	0.00

System Monitoring Compounds

28) TCMX (SURR)	0.00	244	0	0.00	ug/L	
Spiked Amount	10.000		Recovery	=	0.00%	
38) 2,4-DDD (SURR)	27.28	235	122232	10.47	ug/L	0.01
Spiked Amount	10.000		Recovery	=	104.70%	

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	17.47	165	0	N.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.	
29) Azobenzene/ 1,2-Diphenylhyd	0.00	77	0	N.D.	
31) N-Nitrosodiphenylamine	0.00	169	0	N.D.	
32) 4-Bromophenyl-phenyl ether	0.00	248	0	N.D.	
33) Hexachlorobenzene	0.00	284	0	N.D.	
34) Phenanthrene	0.00	178	0	N.D.	
35) Anthracene	0.00	178	0	N.D.	
36) Di-n-butylphthalate	0.00	149	0	N.D.	
37) Fluoranthene	0.00	202	0	N.D.	
40) Pyrene	0.00	202	0	N.D.	
41) Butylbenzylphthalate	0.00	149	0	N.D.	
42) Benzo (a) anthracene	0.00	228	0	N.D. d	
43) Bis (2-ethylhexyl) phthala	30.78	149	2714	0.08 ug/L	95
44) Chrysene	0.00	228	0	N.D. d	

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

6585.D 5080402BBC.M Wed Apr 17 16:44:35 2002

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Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\041502\6585.D

Vial: 9

Acq On : 15 Apr 2002 4:38 pm

Operator: Bohlk

Sample : SAMPLE 6585 3/20/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 17 16:41:14 2002

Quant Results File: 5080402BBC.RES

Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Tue Apr 16 18:27:45 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
46) Di-n-Octylphthalate	0.00	149	0	N.D.		
47) Benzo (b) fluoranthene	0.00	252	0	N.D.		
48) Benzo (k) fluoranthene	0.00	252	0	N.D.		
49) Benzo (a) pyrene	0.00	252	0	N.D.	d	
50) Indeno (1,2,3-cd) pyrene	0.00	276	0	N.D.		
51) Dibenzo (a,h) anthracene	0.00	278	0	N.D.		
52) Benzo (g,h,i) perylene	0.00	276	0	N.D.		

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

6585.D 5080402BBC.M Wed Apr 17 16:44:35 2002

Page 2

Data File : C:\MSDCHEM\1\DATA\041502\6585.D

Vial: 9

Acq On : 15 Apr 2002 4:38 pm

Operator: Bohlk

Sample : SAMPLE 6585 3/20/02

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Apr 17 16:44 2002

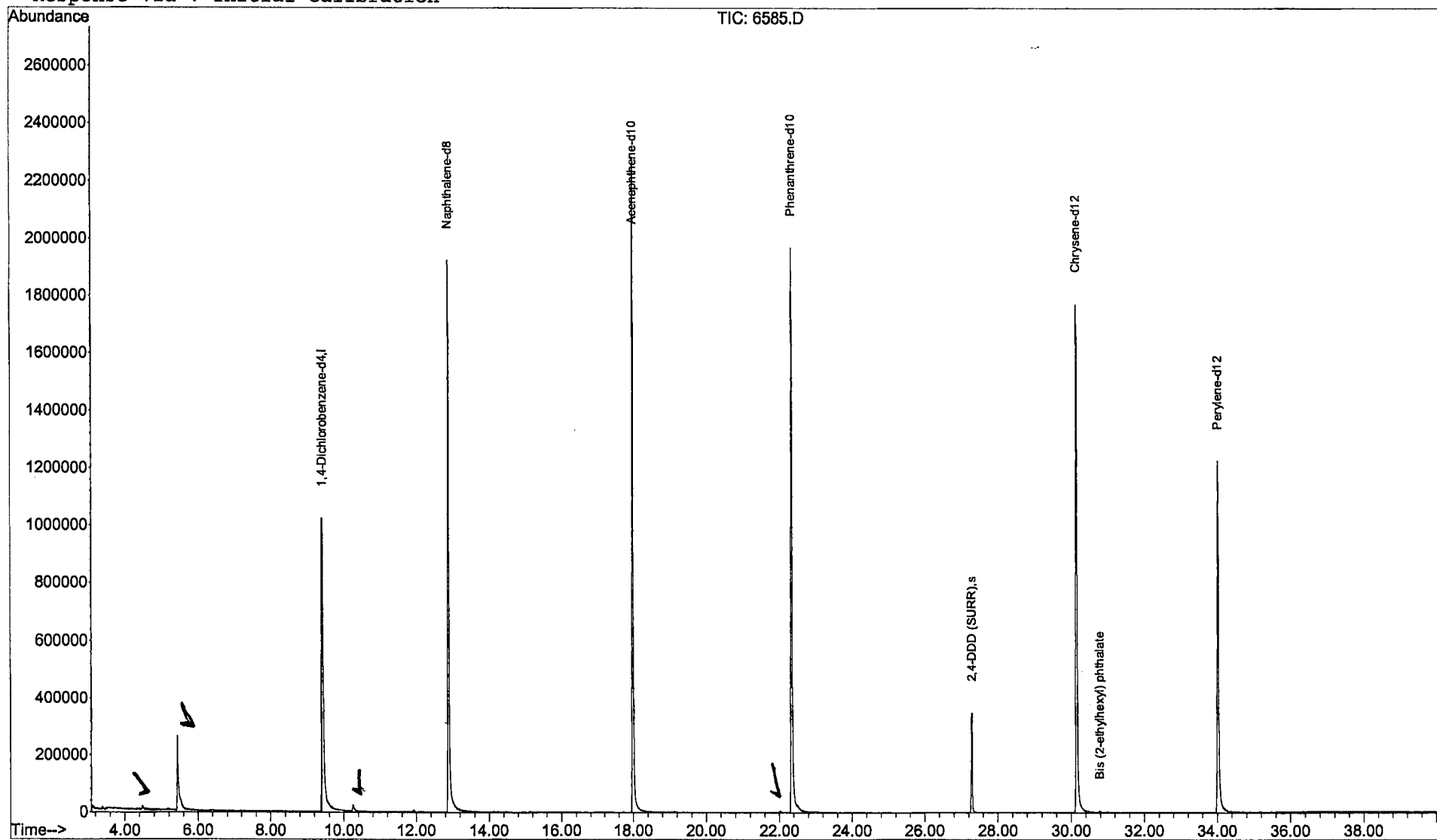
Quant Results File: 5080402BBC.RES

Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Tue Apr 16 18:27:45 2002

Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\041502\6585.D

Acq On : 15 Apr 2002 4:38 pm

Sample : SAMPLE 6585 3/20/02

Misc :

MS Integration Params: LSCINT.P

Vial: 9

Operator: Bohlk

Inst : Instrumen

Multiplr: 1.00

Method : C:\MSDCHEM\1\5973N\5080402BBC.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

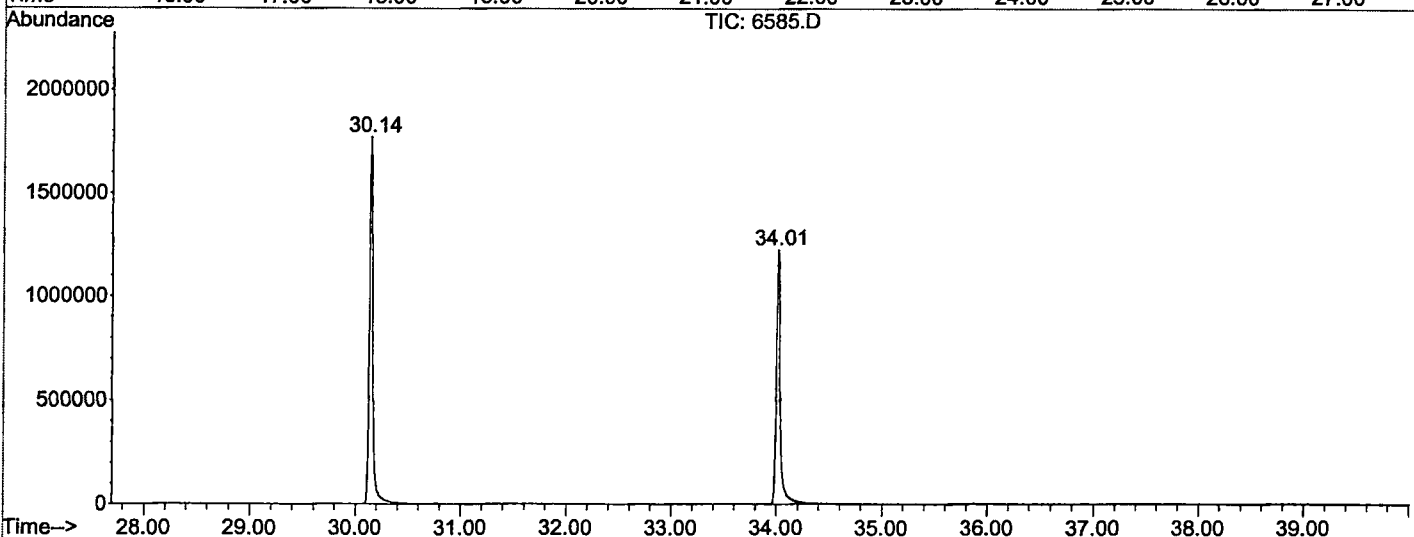
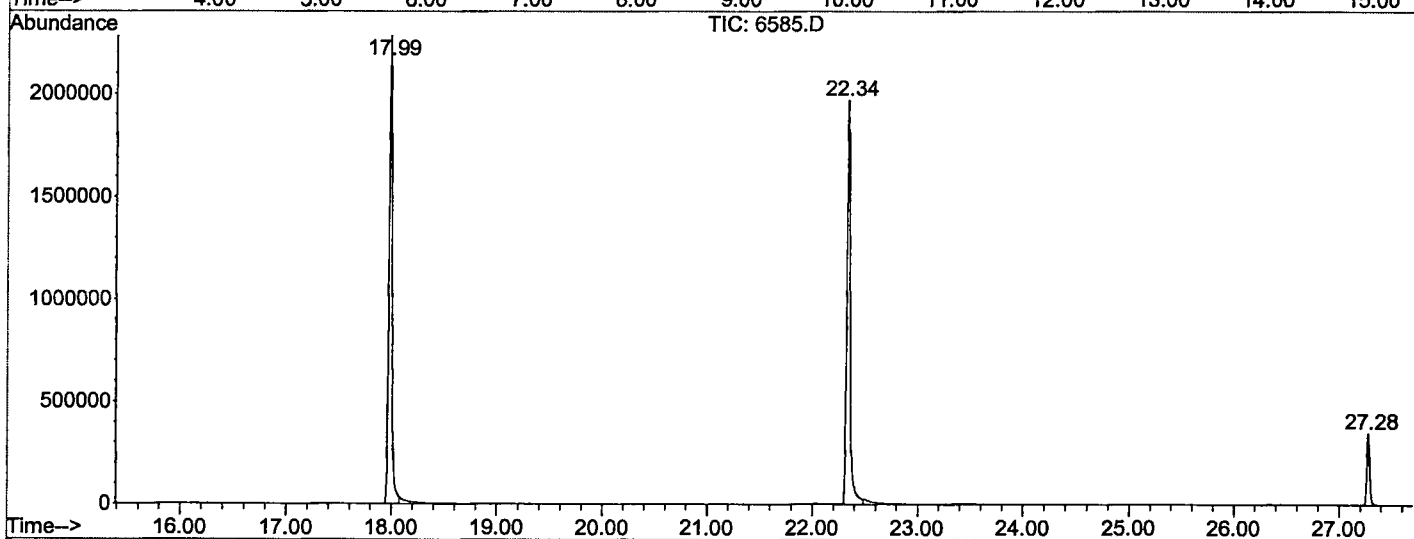
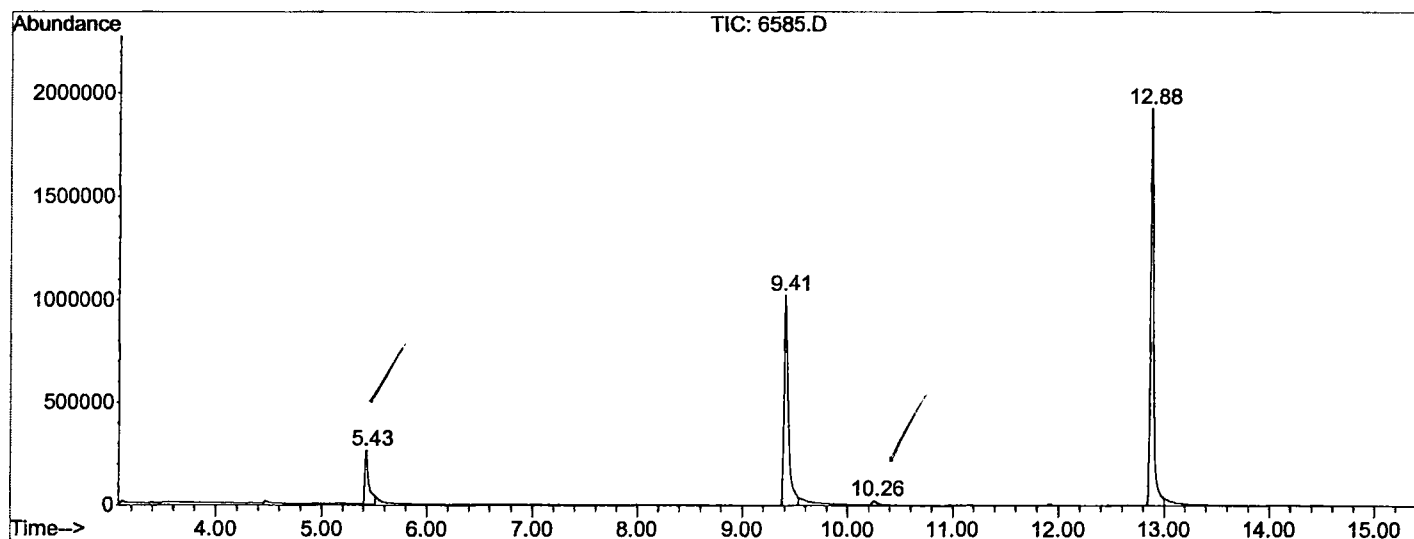
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.429	395	400	414	rBV	260930	647556	14.09%	2.609%
2	9.407	1071	1077	1099	rBV	1021622	2787004	60.64%	11.228%
3	10.259	1213	1222	1231	rBV3	22528	66432	1.45%	0.268%
4	12.879	1661	1668	1689	rBV	1922675	4172122	90.78%	16.809%
5	17.985	2529	2537	2552	rBV2	2277659	4595820	100.00%	18.516%
6	22.339	3269	3278	3303	rBV2	1966805	4524119	98.44%	18.227%
7	27.280	4111	4119	4133	rBV	347690	656973	14.30%	2.647%
8	30.142	4597	4606	4641	rBV2	1769483	4245566	92.38%	17.105%
9	34.014	5255	5265	5294	rBV	1224513	3125471	68.01%	12.592%

Sum of corrected areas: 24821063

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\041502\6585.D
 Operator : Bohlk
 Acquired : 15 Apr 2002 4:38 pm using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: SAMPLE 6585 3/20/02
 Misc Info :
 Vial Number: 9
 Quant File :5080402BBC.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\041502\6585.D

Acq On : 15 Apr 2002 4:38 pm

Sample : SAMPLE 6585 3/20/02

Misc :

MS Integration Params: LSCINT.P

Vial: 9

Operator: Bohlk

Inst : Instrumen

Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)

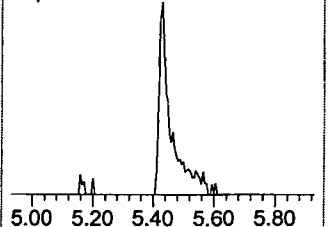
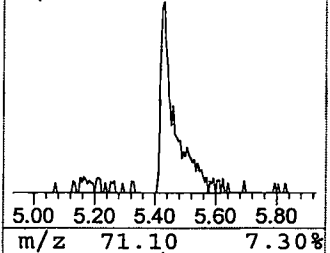
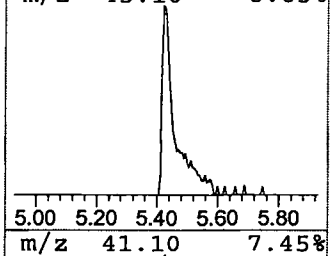
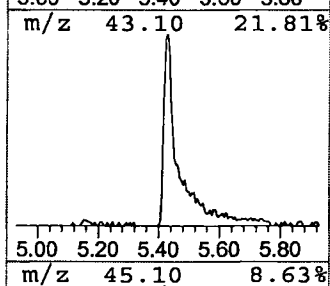
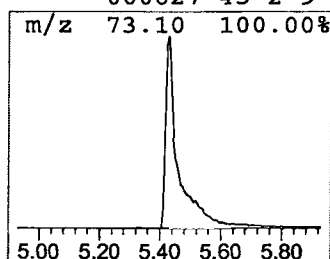
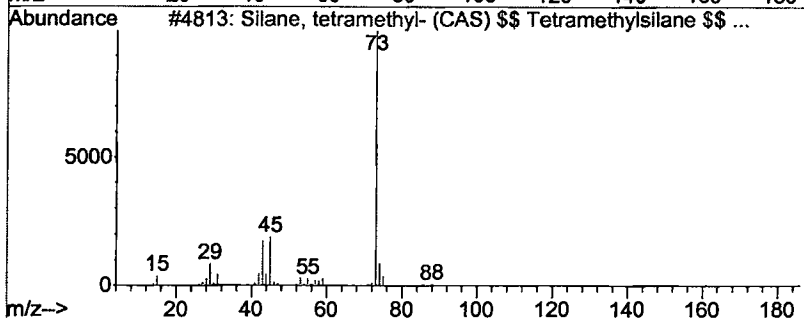
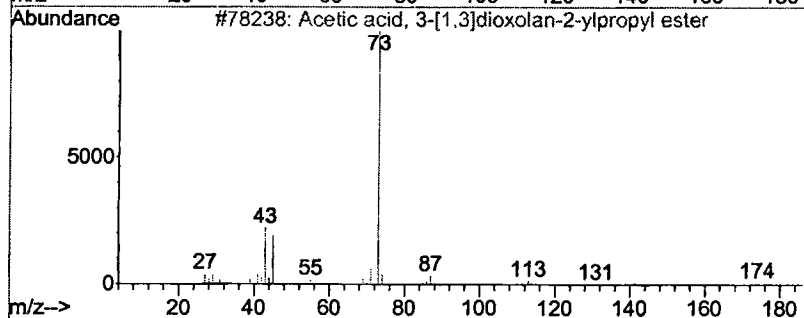
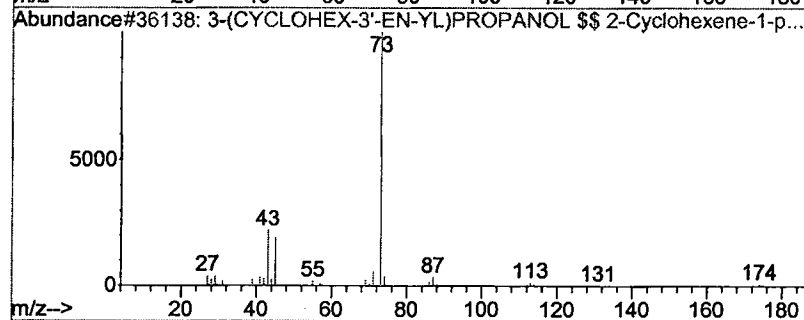
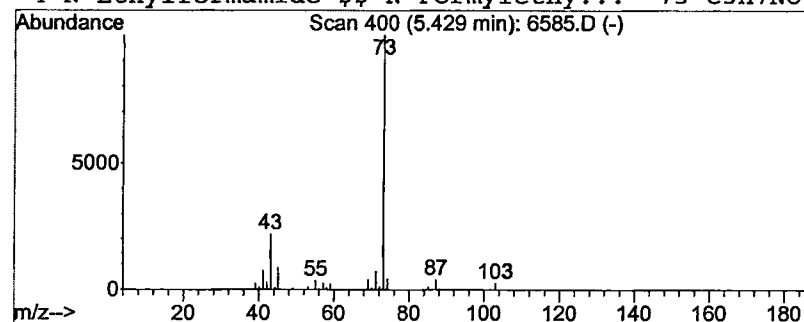
Title : Quantitation For Method 508gcscan

Library : C:\DATABASE\WILEY7N.L

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.43	9.29 ug/L	647556	1,4-Dichlorobenzene-d4	9.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-(CYCLOHEX-3'-EN-YL)PROPANOL \$\$...	174	C8H14O4	015745-87-6	39
2	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	39
3	Silane, tetramethyl- (CAS) \$\$ Te...	88	C4H12Si	000075-76-3	23
4	N-Ethylformamide \$\$ N-Formylethy...	73	C3H7NO	000627-45-2	9



Library Search Compound Report

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Acq On : 15 Apr 2002 4:38 pm
Sample : SAMPLE 6585 3/20/02
Misc :
MS Integration Params: LSCINT.P

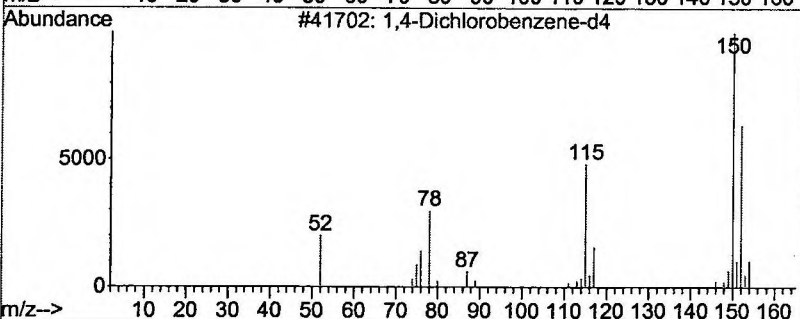
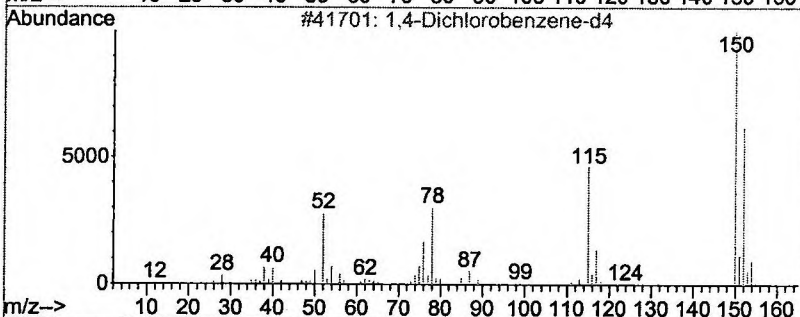
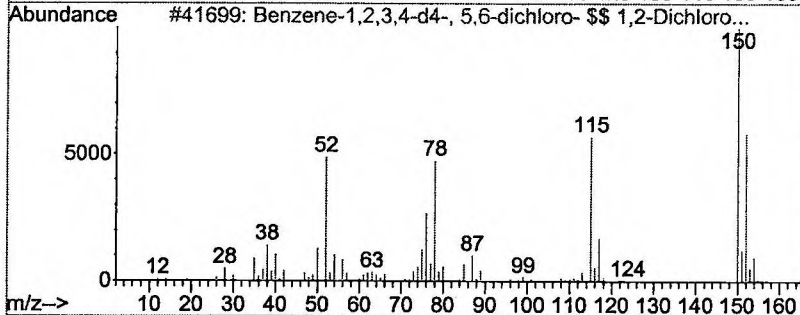
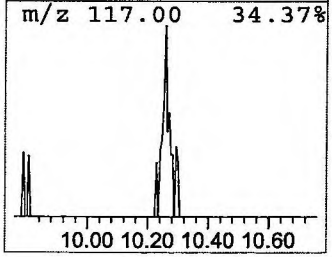
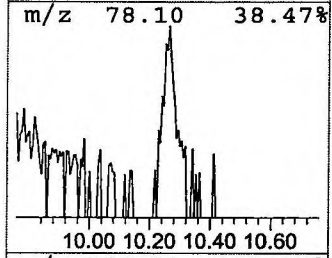
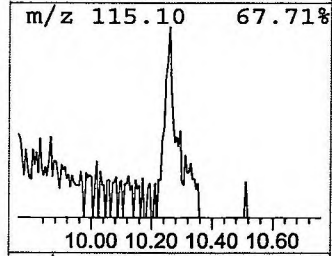
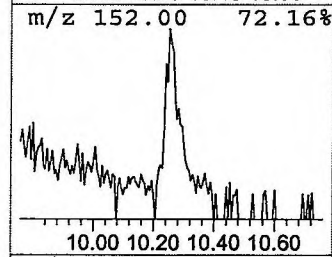
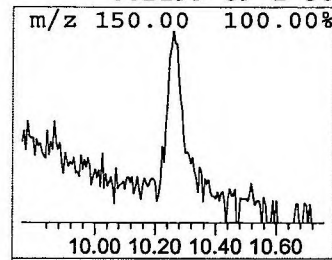
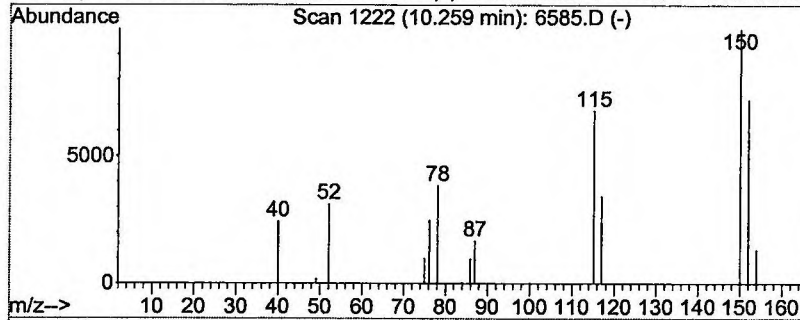
Vial: 9
Operator: Bohlk
Inst : Instrumen
Multiplr: 1.00

Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Library : C:\DATABASE\WILEY7N.L

Peak Number 2 Benzene-1,2,3,4-d4-, 5,6-di... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	0.95 ug/L	66432	1,4-Dichlorobenzene-d4	9.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene-1,2,3,4-d4-, 5,6-dichlor...	150	C6D4Cl2	002199-69-1	91
2	1,4-Dichlorobenzene-d4	150	C6D4Cl2	003855-82-1	91
3	1,4-Dichlorobenzene-d4	150	C6D4Cl2	003855-82-1	86
4	1,2-Dichlorobenzene-d4 \$\$ Benzen...	150	C6D4Cl2	002199-69-1	58



Tentatively Identified Compound (LSC) summary

Operator ID: Bohlk Date Acquired: 15 Apr 2002 4:38 pm
Data File: C:\MSDCHEM\1\DATA\041502\6585.D
Name: SAMPLE 6585 3/20/02
Misc:
Method: C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title: Quantitation For Method 508gcscan
Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response		--Internal Standard---		
					#	RT	Resp	Conc
3-(CYCLOHEX-3'-EN...	5.43	9.3	ug/L	647556	1	9.41	2787000	40.0
Benzene-1,2,3,4-d...	10.26	1.0	ug/L	66432	1	9.41	2787000	40.0