User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/19/2002 Time: 10:34:38

Sample: AE06799
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/19/02 10:34 Ended: 4/19/02 10:34 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		109		5.0

Comments for Sample AE06799 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 04-19-2002 Time: 10:35:03

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\041602\6799.D Vial: 17
Acq On : 16 Apr 2002 9:48 pm Operator: Bohlk
Sample : SAMPLE 6799 3/21/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 18 18:09:58 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	412441	40.00	ug/L	0.00
10) Naphthalene-d8	12.88	136	1799877	40.00	ug/L	0.00
17) Acenaphthene-d10	17.99	164	1045220	40.00	ug/L	0.00
30) Phenanthrene-d10	22.33	188	1528097	40.00	ug/L	0.00
39) Chrysene-d12	30.14	240	1274627	40.00	ug/L	-0.01
45) Perylene-d12	34.01	264	847616	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	123788	10.88	ug/L	0.01
Spiked Amount 10.000			Recovery = 108.80%			
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.		
29) Azobenzen/ 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.	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) antracene	0.00	228	0	N.D.	d	
43) Bis (2-ethylhexyl) phthala	30.76	149	5752	0.18	ug/L	87
44) Chrysene	0.00	228	0	N.D.	d	

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

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Sample : SAMPLE 6799 3/21/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 18 18:09:58 2002 Quant Results File: 5080402BBC.RES

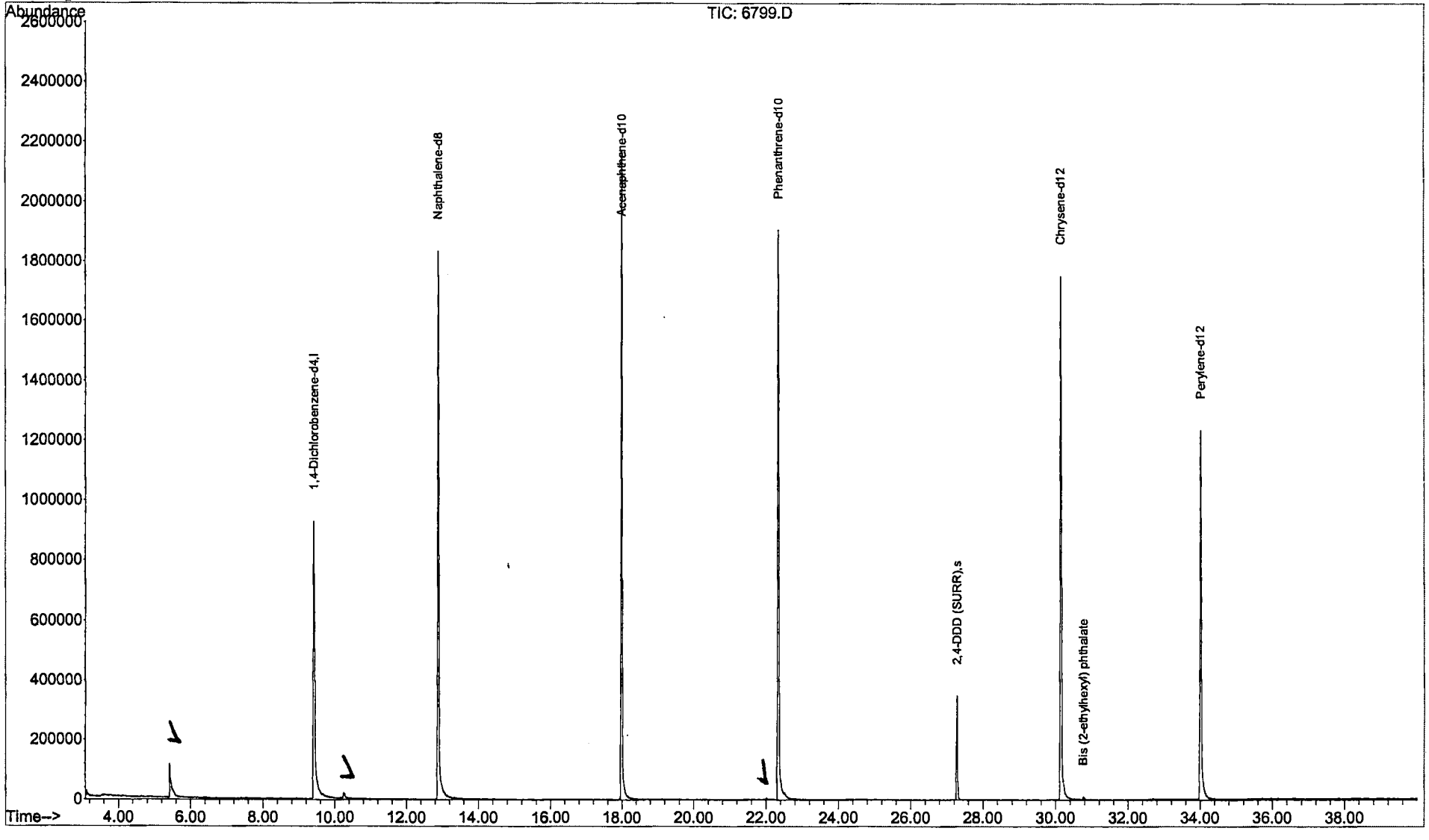
Quant Method : C:\MSDCHEM\1\5973N\5080402BBC.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Tue Apr 16 18:27:45 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

	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)	Dibenz (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

Data File : C:\MSDCHEM\1\DATA\041602\6799.D Vial: 17
Acq On : 16 Apr 2002 9:48 pm Operator: Bohlk
Sample : SAMPLE 6799 3/21/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Apr 18 18:11 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



Data File : C:\MSDCHEM\1\DATA\041602\6799.D Vial: 17
Acq On : 16 Apr 2002 9:48 pm Operator: Bohlk
Sample : SAMPLE 6799 3/21/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

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

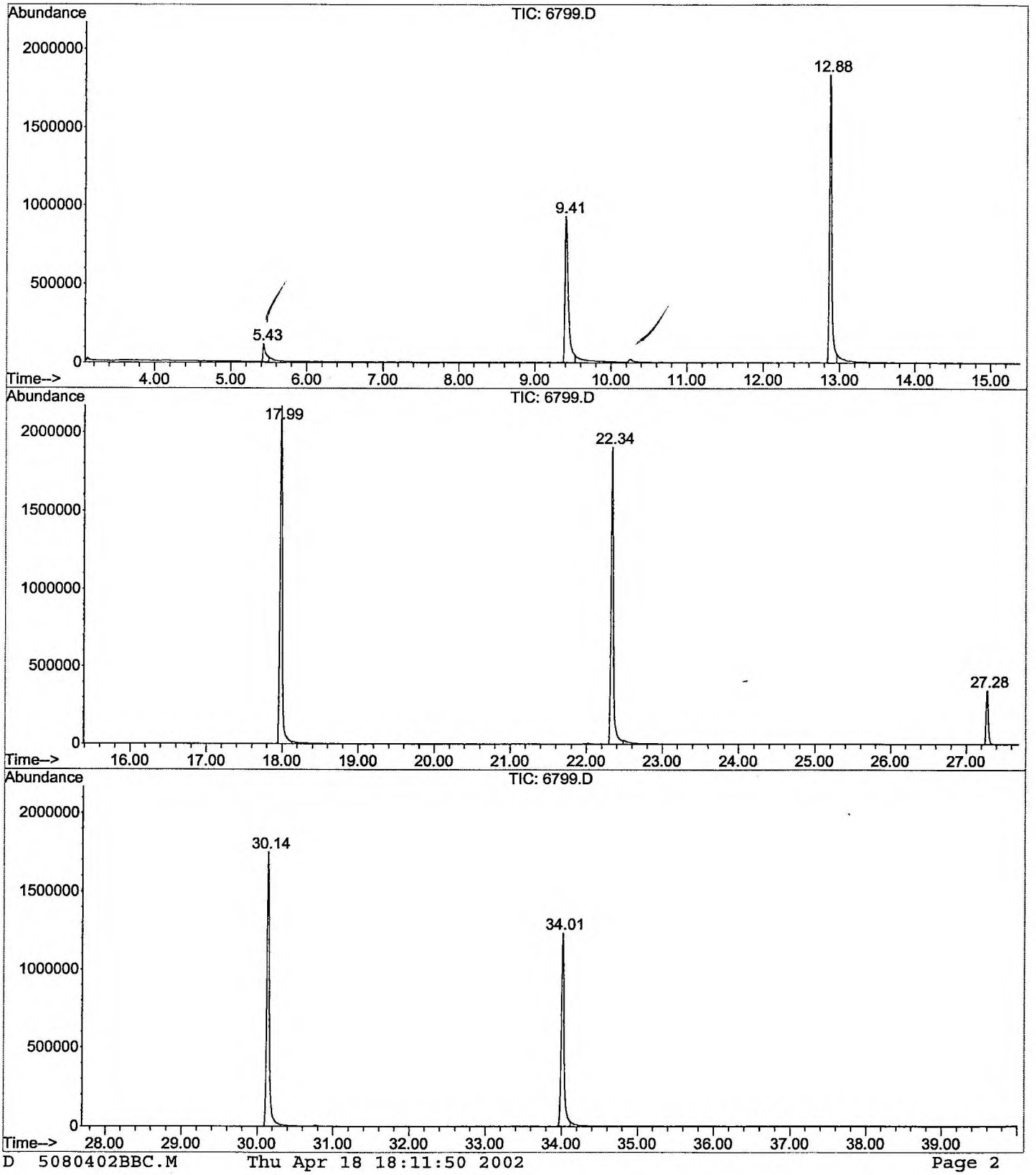
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.429	396	400	412	rBV	111695	296218	6.57%	1.256%
2	9.407	1071	1077	1097	rBV	925762	2627355	58.26%	11.144%
3	12.879	1661	1668	1683	rBV	1834084	3890280	86.26%	16.501%
4	17.985	2529	2537	2562	rBV	2167911	4510073	100.00%	19.129%
5	22.339	3269	3278	3302	rBV2	1903803	4377176	97.05%	18.566%
6	27.280	4111	4119	4135	rBV2	347121	666506	14.78%	2.827%
7	30.142	4596	4606	4635	rBV2	1749896	4127173	91.51%	17.505%
8	34.014	5256	5265	5283	rBV2	1234913	3081842	68.33%	13.072%

Sum of corrected areas: 23576623

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\041602\6799.D
Operator : Bohlk
Acquired : 16 Apr 2002 9:48 pm using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: SAMPLE 6799 3/21/02
Misc Info :
Vial Number: 17
Quant File :5080402BBC.RES (RTE Integrator)



Data File : C:\MSDCHEM\1\DATA\041602\6799.D Vial: 17
Acq On : 16 Apr 2002 9:48 pm Operator: Bohlk
Sample : SAMPLE 6799 3/21/02 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

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

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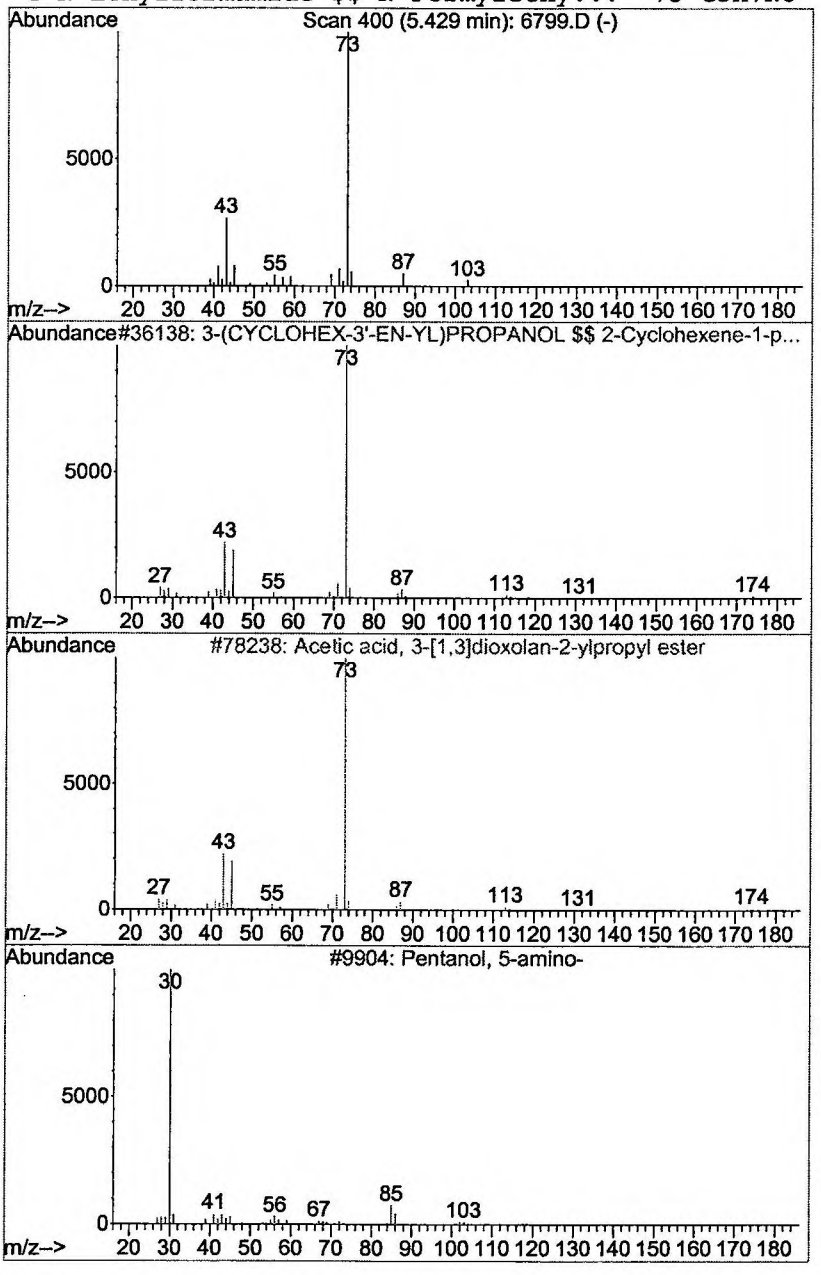
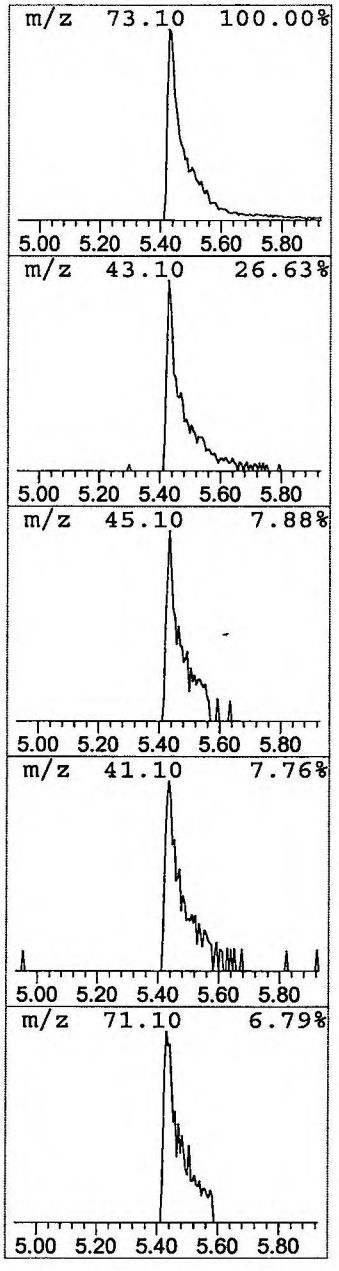
Peak Number 1 3-(CYCLOHEX-3'-EN-YL)PROPAN... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.43	4.51 ug/L	296218	1,4-Dichlorobenzene-d4	9.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-(CYCLOHEX-3'-EN-YL)PROPANOL $$...	174	C8H14O4	015745-87-6	36
2	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	36
3	Pentanol, 5-amino-	103	C5H13NO	002508-29-4	9
4	N-Ethylformamide $$ N-Formylethy...	73	C3H7NO	000627-45-2	9





Tentatively Identified Compound (TIC) Summary

Operator ID: Bohlk Date Acquired: 16 Apr 2002 9:48 pm
Data File: C:\MSDCHEM\1\DATA\041602\6799.D
Name: SAMPLE 6799 3/21/02
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 #	Internal Standard RT	Internal Standard Resp	Internal Standard Conc
3-(CYCLOHEX-3'-EN...	5.43	4.5	ug/L	296218	1	9.41	2627360	40.0