

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/22/2002 Time: 17:43:31

Sample: AE06267
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/22/02 17:32 Ended: 3/22/02 17:32 Analyst: DLV
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD	USPEC	151		5.0

Comments for Sample AE06267 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-22-2002 Time: 17:43:33

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for the LCS Standard compounds are above their acceptance ranges. New control limits will be calculated.

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031602\06267.D

Acq On : 16 Mar 2002 12:31 pm

Sample : 3/15

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 20 06:27:09 2002

Vial: 5

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

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	1403788	20.00	ug/L	0.00
10) Naphthalene-d8	13.03	136	3891995	20.00	ug/L	0.00
17) Acenaphthene-d10	18.13	164	2141746	20.00	ug/L	0.00
29) Phenanthrene-d10	22.50	188	3299911	20.00	ug/L	0.00
38) Chrysene-d12	30.30	240	3106014	20.00	ug/L	-0.02
44) Perylene-d12	34.18	264	1726685	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-ddd surr	27.43	235	297074	7.55	ug/L	0.00
Spiked Amount	5.000		Recovery	151.00%		

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.	
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) anthracene	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

06267.D 5080302.M Wed Mar 20 06:29:48 2002

Quantitation Report (QT Reviewed)

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Misc :

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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

06267.D 5080302.M Wed Mar 20 06:29:49 2002

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\031602\06267.D

Vial: 5

Acq On : 16 Mar 2002 12:31 pm

Operator: McLean

Sample : 3/15

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Mar 20 6:29 2002

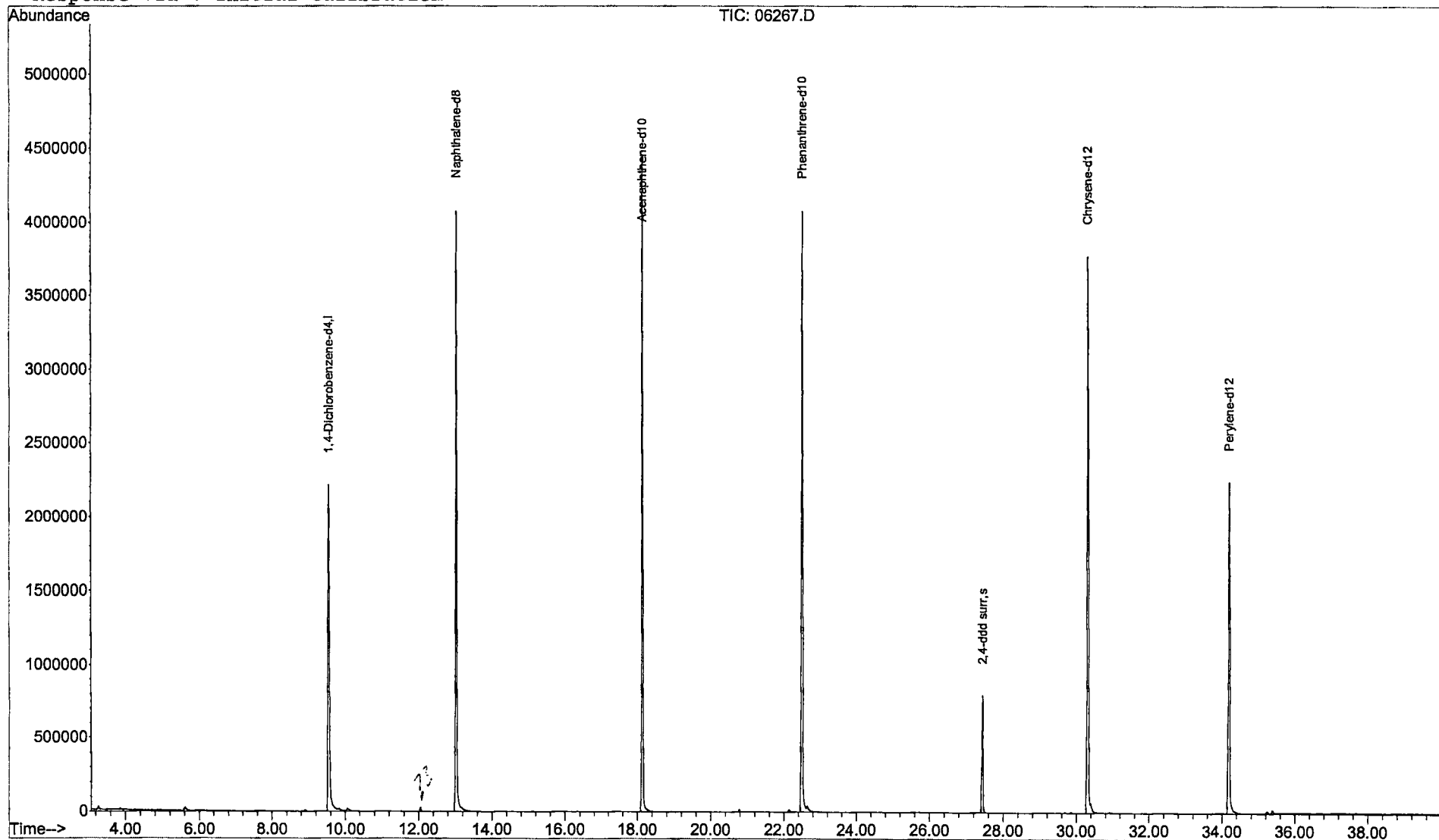
Quant Results File: 5080302.RES

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\031602\06267.D

Acq On : 16 Mar 2002 12:31 pm

Sample : 3/15

Misc :

MS Integration Params: LSCINT.P

Vial: 5

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

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

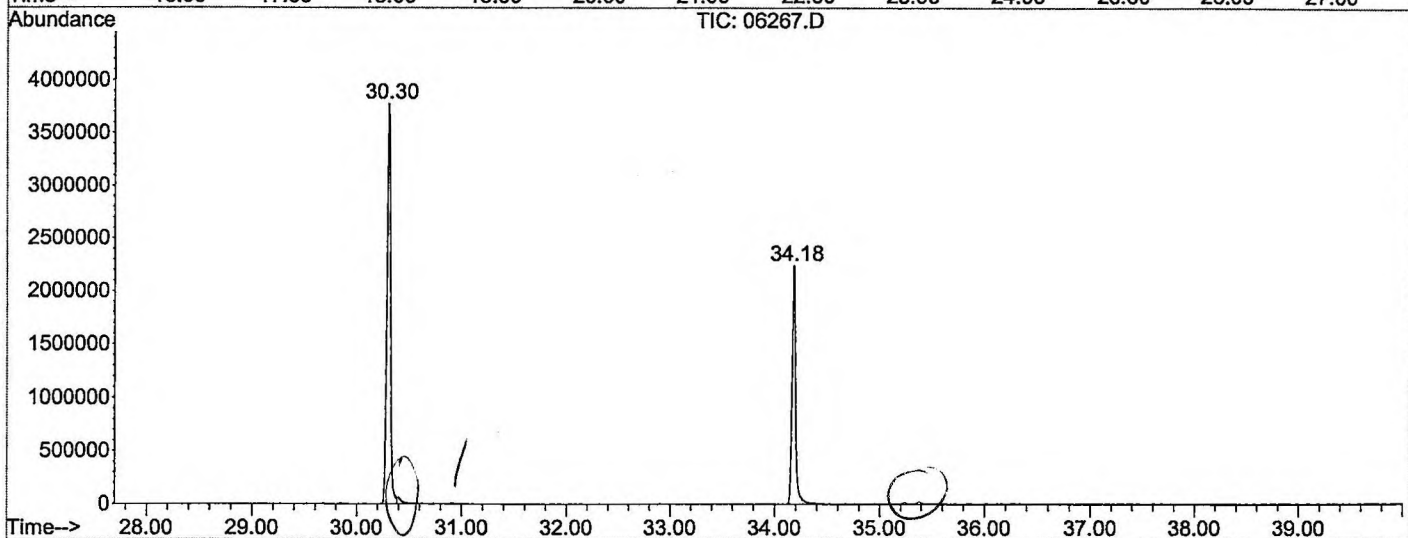
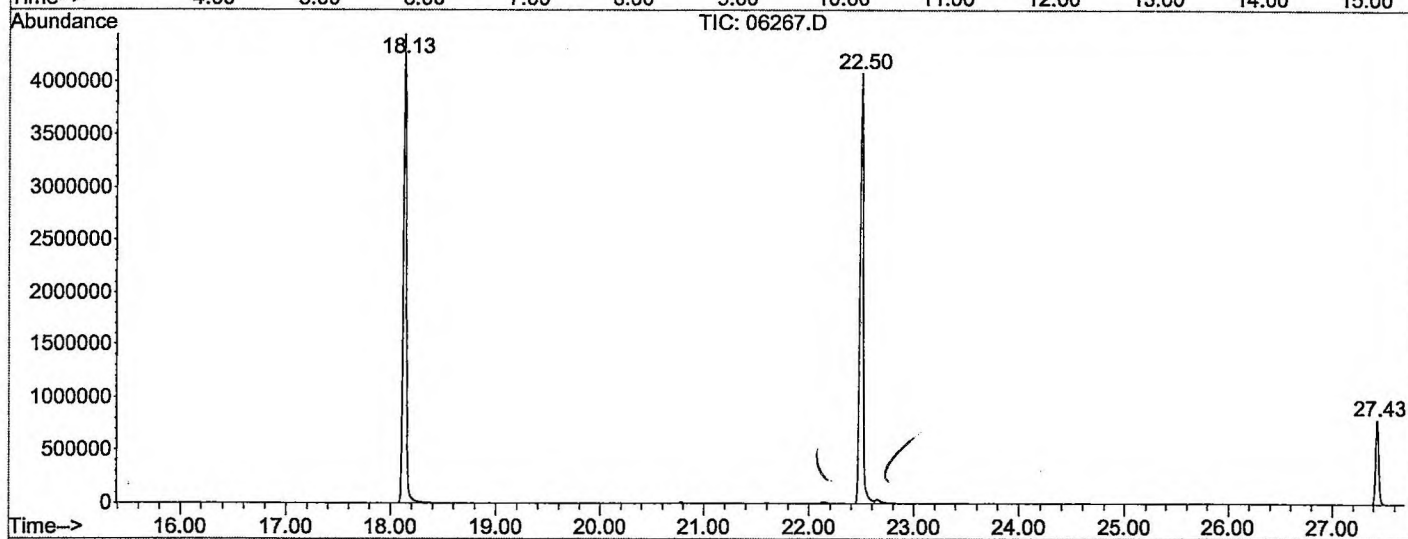
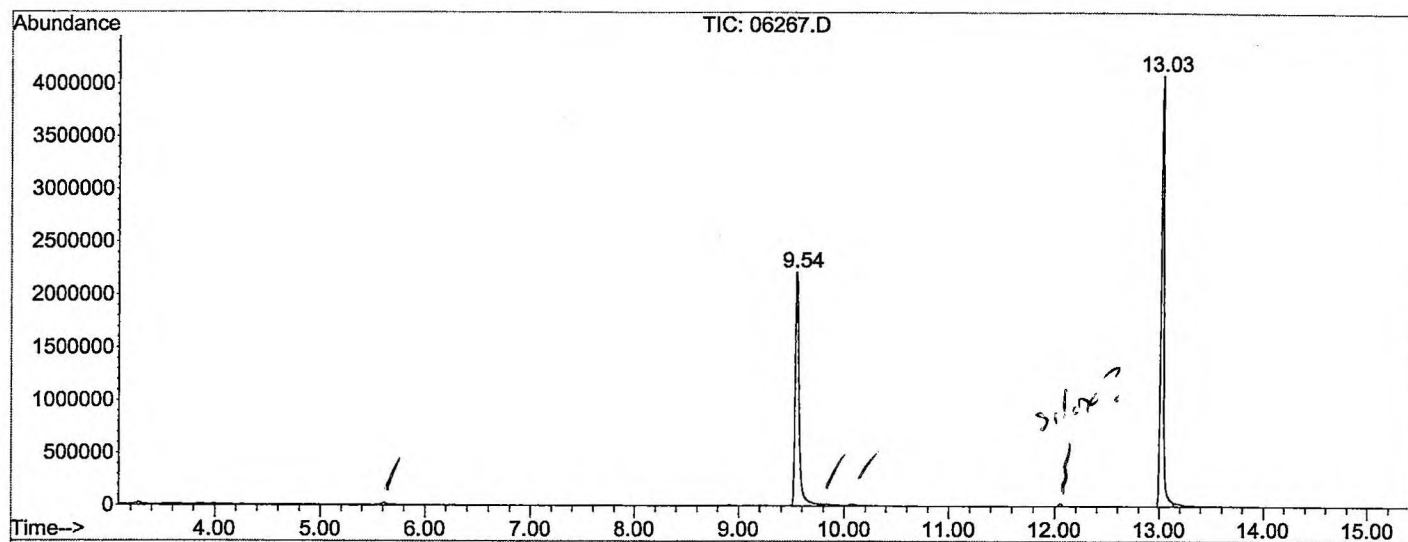
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.543	708	713	736	rBV	2211932	5953471	65.56%	12.347%
2	13.026	1091	1097	1109	rBV	4083413	8139932	89.64%	16.881%
3	18.133	1654	1660	1679	rBV	4449286	8898261	97.99%	18.454%
4	22.495	2135	2141	2155	rBV2	4079389	9080838	100.00%	18.833%
5	27.430	2680	2685	2698	rBB	793660	1542754	16.99%	3.200%
6	30.305	2995	3002	3011	rBV2	3774993	8953424	98.60%	18.569%
7	34.178	3423	3429	3444	rBV2	2241451	5649593	62.21%	11.717%

Sum of corrected areas: 48218273

LSC Report - Integrated Chromatogram

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File       : C:\MSDCHEM\1\DATA\031602\06267.D
Operator   : McLean
Acquired    : 16 Mar 2002  12:31 pm using AcqMethod 508SCAN
Instrument  : Instrumen
Sample Name: 3/15
Misc Info  :
Vial Number: 5
Quant File : 5080302.RES (RTE Integrator)
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Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 16 Mar 2002 12:31 pm
 Data File: C:\MSDCHEM\1\DATA\031602\06267.D
 Name: 3/15
 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	#	RT	Resp	Conc
