

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/04/2002 Time: 13:53:32

Sample: AE01641
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/4/02 13:53 Ended: 3/4/02 13:53 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		76		5.0

Comments for Sample AE01641 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 03-04-2002 Time: 13:53:58

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 10 of the 43 compounds in the LCS Standard were below their acceptance ranges.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\022602\1641.D

Vial: 7

Acq On : 26 Feb 2002 2:58 pm

Operator: McLean

Sample : 1/24

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Feb 28 11:24:08 2002

Quant Results File: 5080202.RES

Quant Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Feb 25 19:17:14 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	1197949	20.00	ug/L	0.00
10) Naphthalene-d8	13.05	136	3181046	20.00	ug/L	0.00
17) Acenaphthene-d10	18.17	164	1720048	20.00	ug/L	0.00
29) Phenanthrene-d10	22.52	188	2653224	20.00	ug/L	0.00
38) Chrysene-d12	30.34	240	2545040	20.00	ug/L	-0.02
44) Perylene-d12	34.21	264	1733972	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-ddd surr	27.47	235	159619	3.79	ug/L	0.00
Spiked Amount	5.000		Recovery	=	75.80%	

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	24.67	149	9784	0.06 ug/L	79
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.	
43) Chrysene	30.34	228	6182	0.04 ug/L	57
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

1641.D 5080202.M Thu Feb 28 11:24:36 2002

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\022602\1641.D

Vial: 7

Acq On : 26 Feb 2002 2:58 pm

Operator: McLean

Sample : 1/24

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Feb 28 11:24:08 2002

Quant Results File: 5080202.RES

Quant Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)

Title : Quantitation For Method 508gcscan

Last Update : Mon Feb 25 19:17:14 2002

Response via : Initial Calibration

DataAcq Meth : 508SCAN

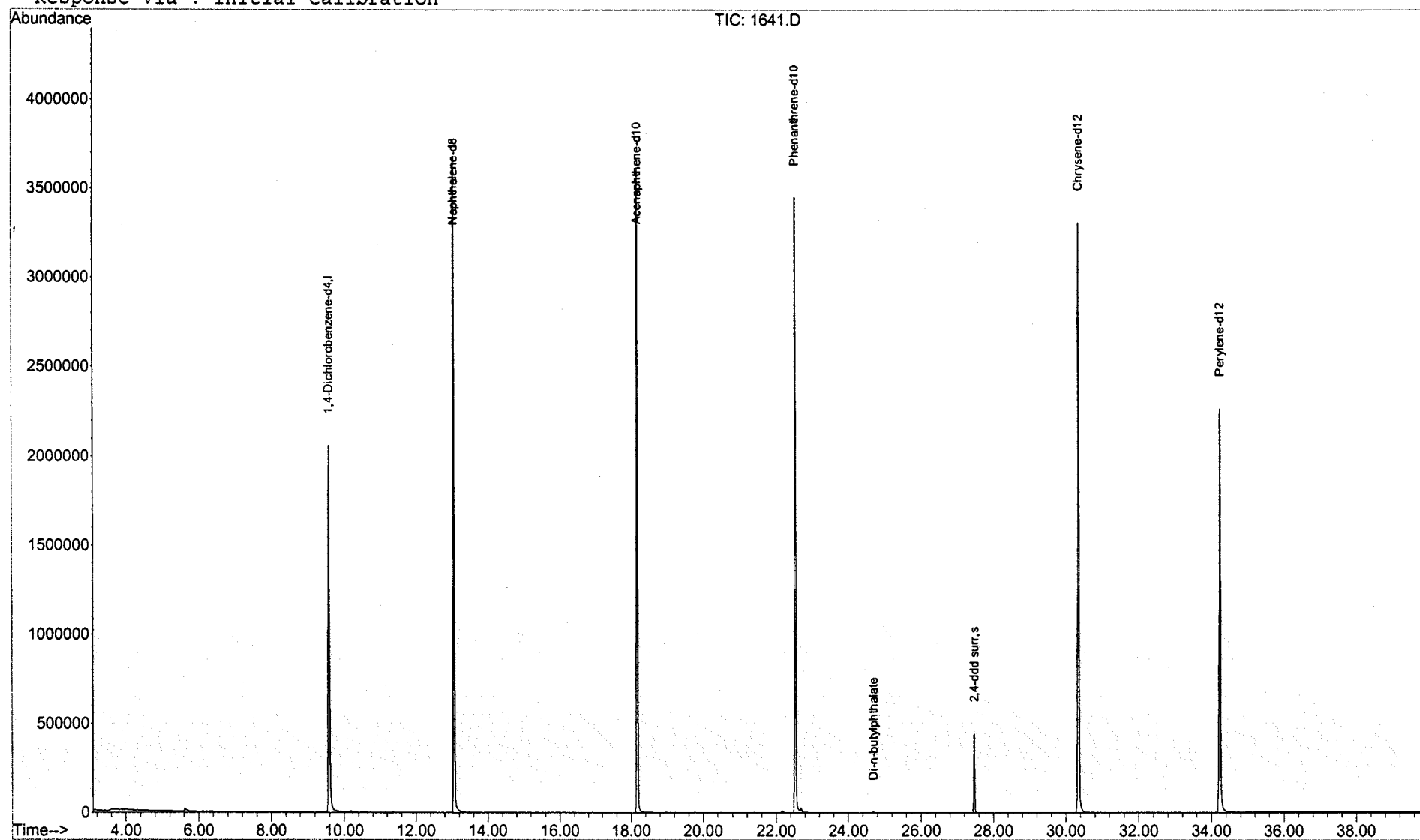
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
47) Benzo (k) fluoranthene	0.00	252	0	N.D.		
48) Benzo (a) pyrene	0.00	252	0	N.D.	d	
49) Indeno (1,2,3-cd) pyrene	0.00	276	0	N.D.		
50) Dibenz (a,h) anthracene	0.00	278	0	N.D.		
51) Benzo (g,h,i) perylene	0.00	276	0	N.D.		

(#) = qualifier out of range (m) = manual integration
1641.D 5080202.M Thu Feb 28 11:24:36 2002

Page 2

Data File : C:\MSDCHEM\1\DATA\022602\1641.D Vial: 7
Acq On : 26 Feb 2002 2:58 pm Operator: McLean
Sample : 1/24 Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Feb 28 11:24 2002 Quant Results File: 5080202.RES

Method : C:\MSDCHEM\1\5973N\5080202.M (RTE Integrator)
Title : Quantitation For Method 508gcscan
Last Update : Mon Feb 25 19:17:14 2002
Response via : Initial Calibration



1641.D 5080202.M Thu Feb 28 11:24:37 2002

Page 3

LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\022602\1641.D
 Acq On : 26 Feb 2002 2:58 pm
 Sample : 1/24
 Misc :
 MS Integration Params: LSCINT.P

Vial: 7
 Operator: McLean
 Inst : Instrumen
 Multiplr: 1.00

Method : C:\MSDCHEM\1\5973N\5080202.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

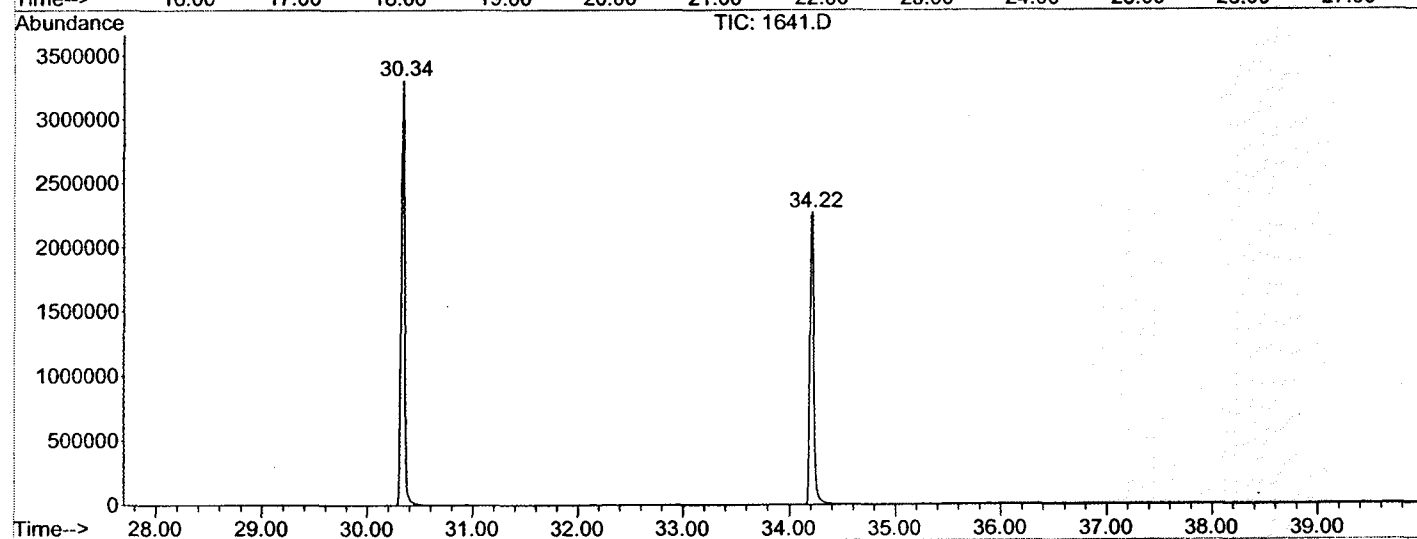
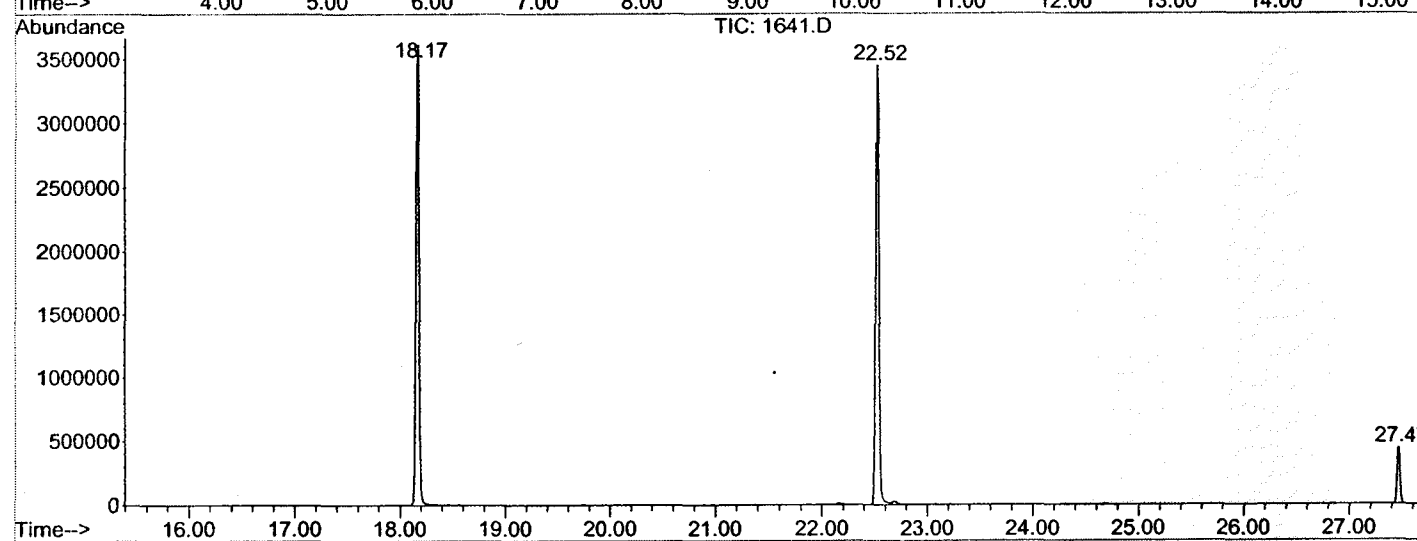
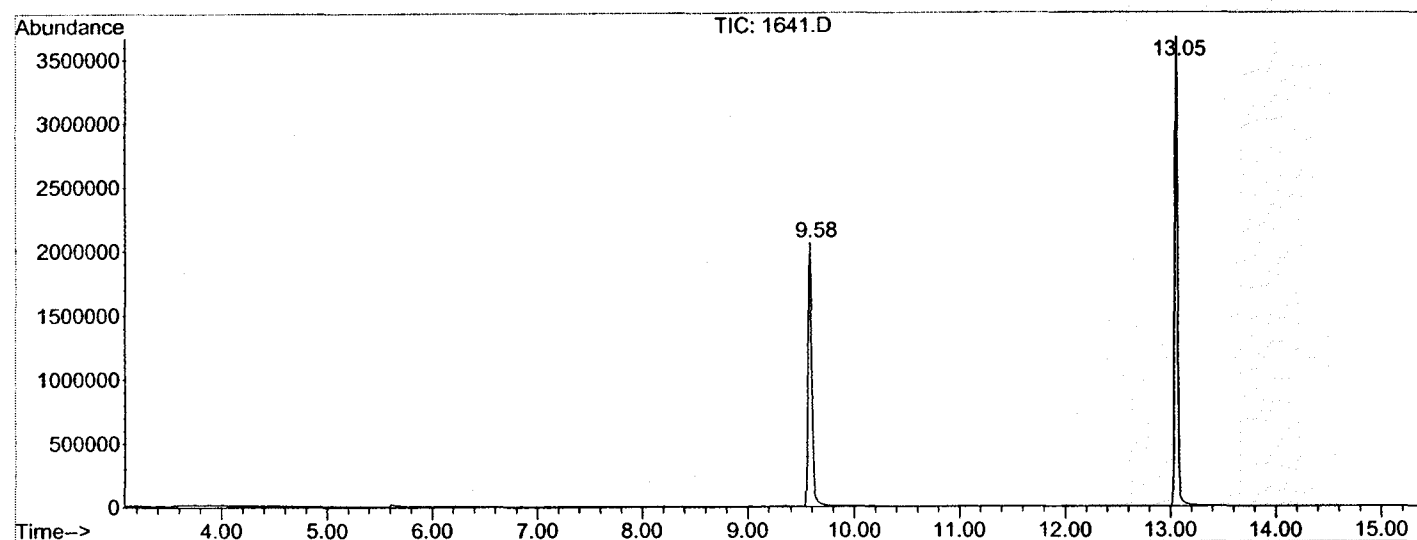
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.579	712	717	735	rBV	2058251	5049504	68.47%	12.489%
2	13.053	1095	1100	1116	rBV	3665686	6882430	93.32%	17.023%
3	18.169	1658	1664	1675	rBV	3615254	7200252	97.63%	17.809%
4	22.523	2138	2144	2157	rBV2	3449859	7331644	99.41%	18.134%
5	27.466	2684	2689	2698	rBB	440409	827595	11.22%	2.047%
6	30.341	2999	3006	3019	rBV2	3306861	7374939	100.00%	18.241%
7	34.223	3426	3434	3450	rBV2	2267151	5763789	78.15%	14.256%

Sum of corrected areas: 40430153

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\022602\1641.D
 Operator : McLean
 Acquired : 26 Feb 2002 2:58 pm using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: 1/24
 Misc Info :
 Vial Number: 7
 Quant File :5080202.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 26 Feb 2002 2:58 pm
 Data File: C:\MSDCHEM\1\DATA\022602\1641.D
 Name: 1/24
 Misc:
 Method: C:\MSDCHEM\1\5973N\5080202.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