

User: Vinci, David L
 Westchester Dept. of Labs & Research
 Date: 03/07/2002 Time: 16:42:59

Sample: AE02566
 Analysis: \$GCMS GCMS Scan For Unknowns
 Units: ug
 Started: 3/7/02 16:42 Ended: 3/7/02 16:42 Analyst: DLV
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds				
2	Surr_2,4-DDD	70			5.0

Comments for Sample AE02566 - Analysis \$GCMS

Newestchester Dept. of Labs & Research
User: Vinci, David L
Date: 03-07-2002 Time: 16:48:06

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 9 of the 43 compounds in the LCS Standard were below their acceptance ranges.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\022702\02566.D

Acq On : 27 Feb 2002 7:45 pm

Sample : GC SCAN 2/4

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 01 08:59:14 2002

Vial: 14

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

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	1642483	20.00	ug/L	0.00
10) Naphthalene-d8	13.05	136	4298150	20.00	ug/L	0.00
17) Acenaphthene-d10	18.17	164	2337273	20.00	ug/L	0.00
29) Phenanthrene-d10	22.53	188	3640858	20.00	ug/L	0.00
38) Chrysene-d12	30.35	240	3489630	20.00	ug/L	0.00
44) Perylene-d12	34.22	264	2470158	20.00	ug/L	0.00

System Monitoring Compounds

37) 2,4-ddd surr	27.47	235	202596	3.50	ug/L	0.00
Spiked Amount	5.000		Recovery	=	70.00%	

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	18.17	165	304144	9.35 ug/L #	25
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	15875	0.07 ug/L	94
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	30.94	149	50555	0.32 ug/L	99
43) Chrysene	30.35	228	8323	0.04 ug/L	46
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

02566.D 5080202.M Fri Mar 01 09:00:02 2002

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\022702\02566.D
 Acq On : 27 Feb 2002 7:45 pm
 Sample : GC SCAN 2/4
 Misc :
 MS Integration Params: BENZO.P
 Quant Time: Mar 01 08:59:14 2002

Vial: 14
 Operator: McLean
 Inst : Instrumen
 Multiplr: 1.00
 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
 02566.D 5080202.M Fri Mar 01 09:00:02 2002

Data File : C:\MSDCHEM\1\DATA\022702\02566.D

Acq On : 27 Feb 2002 7:45 pm

Sample : GC SCAN 2/4

Misc :

MS Integration Params: BENZO.P

Quant Time: Mar 1 8:59 2002

Vial: 14

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

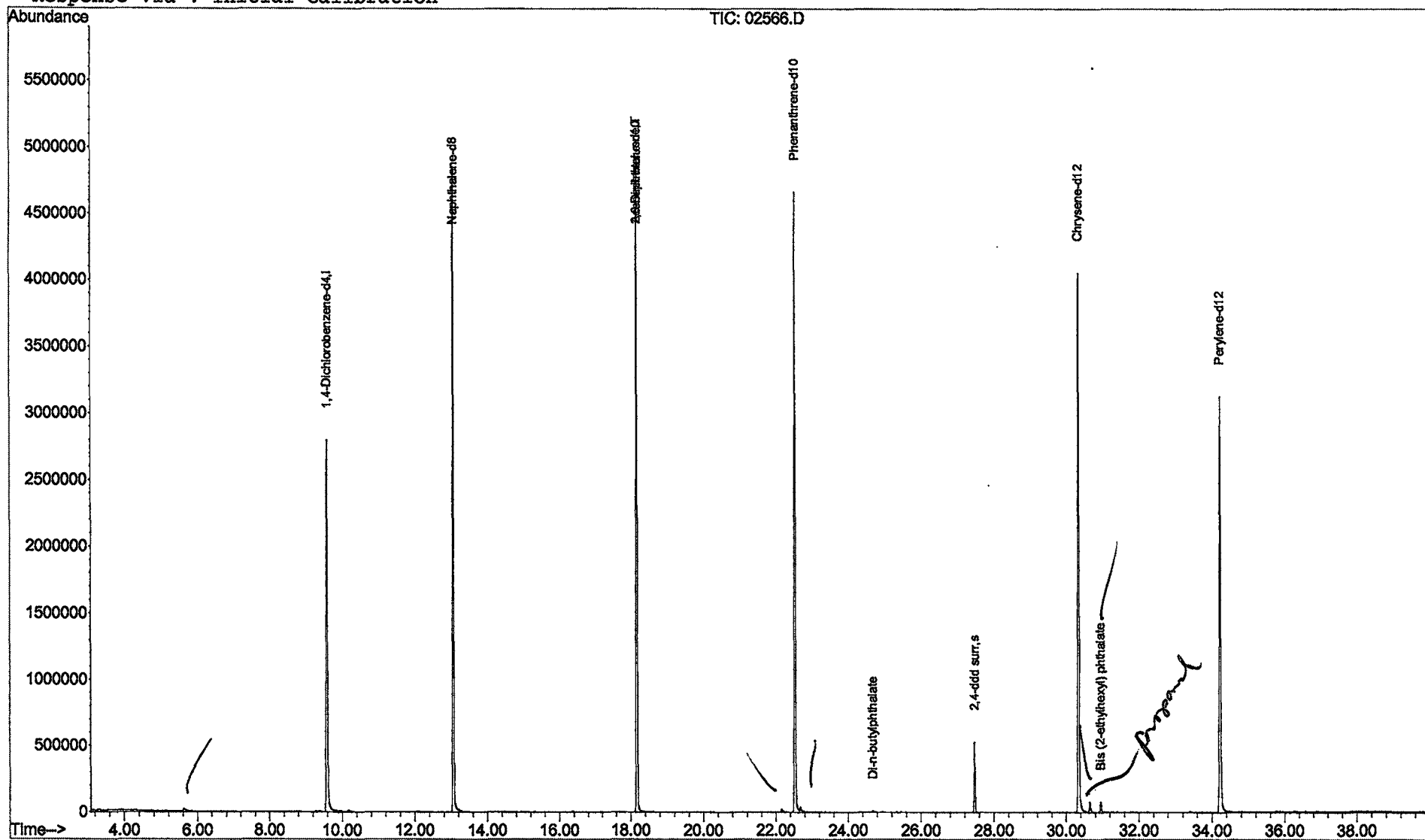
Quant Results File: 5080202.RES

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



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\022702\02566.D Vial: 14
 Acq On : 27 Feb 2002 7:45 pm Operator: McLean
 Sample : GC SCAN 2/4 Inst : Instrumen
 Misc : Multiplr: 1.00
 MS Integration Params: LSCINT.P

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

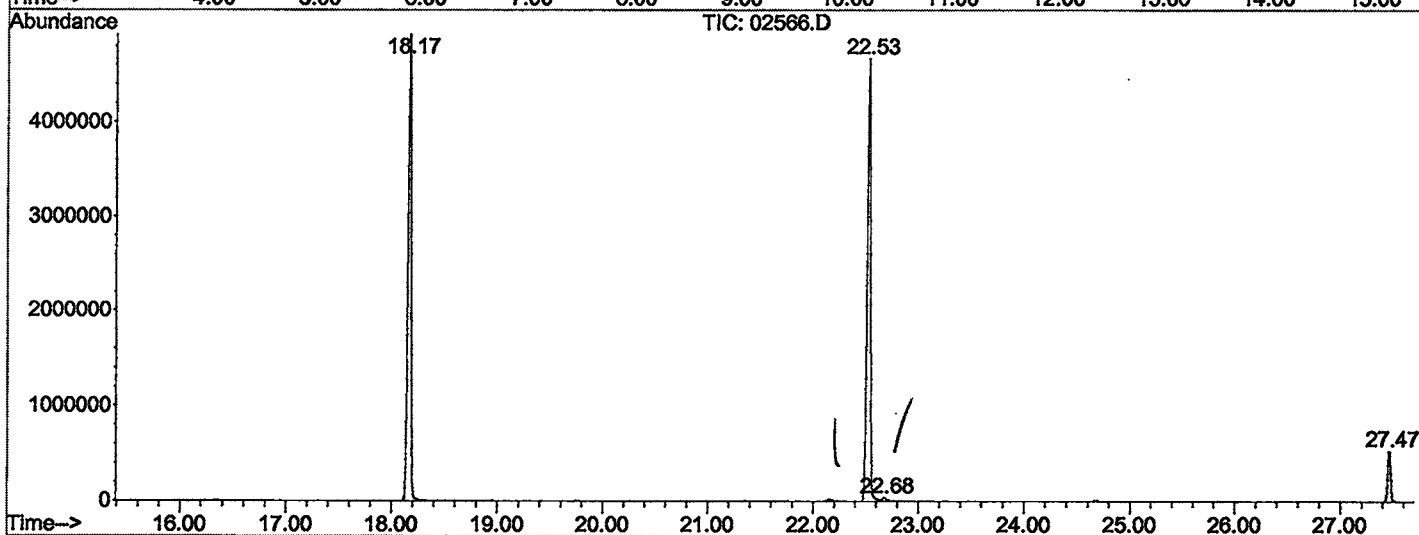
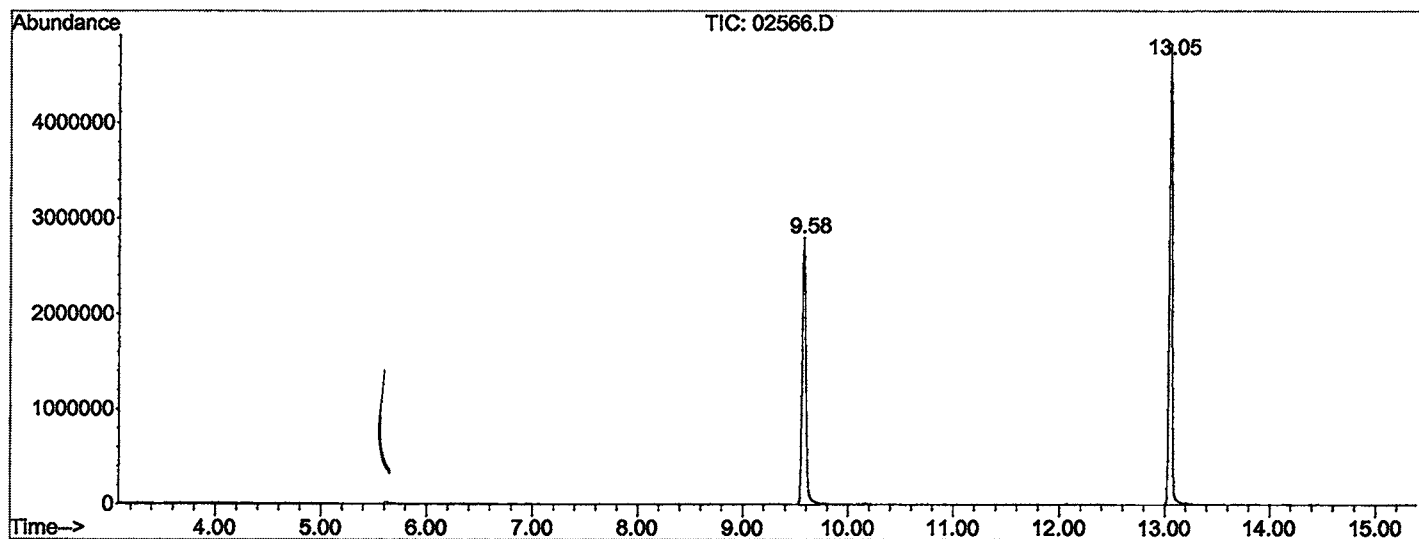
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	9.579	711	717	742	rBV	2796476	6933733	68.16%	12.441%
2	13.053	1095	1100	1113	rBV	4830495	9228264	90.71%	16.558%
3	18.169	1658	1664	1680	rBV	4918147	9762169	95.96%	17.516%
4	22.532	2138	2145	2157	rBV2	4666080	10018609	98.48%	17.976%
5	22.677	2158	2161	2178	rVB	37407	106921	1.05%	0.192%
6	27.466	2684	2689	2699	rBV	529249	1035332	10.18%	1.858%
7	30.350	2999	3007	3025	rBV	4054233	10173363	100.00%	18.254%
8	30.641	3035	3039	3047	rBV	76078	180442	1.77%	0.324%
9	30.940	3066	3072	3082	rVB	77347	179336	1.76%	0.322%
10	34.223	3426	3434	3448	rBV2	3124874	8114644	79.76%	14.560%

Sum of corrected areas: 55732813

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\022702\02566.D
 Operator : McLean
 Acquired : 27 Feb 2002 7:45 pm using AcqMethod 508SCAN
 Instrument : Instrumen
 Sample Name: GC SCAN 2/4
 Misc Info :
 Vial Number: 14
 Quant File :5080202.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: McLean Date Acquired: 27 Feb 2002 7:45 pm
 Data File: C:\MSDCHEM\1\DATA\022702\02566.D
 Name: GC SCAN 2/4
 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