

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 04/22/2002 Time: 09:00:48

Sample: AE06986
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 4/22/02 09:00 Ended: 4/22/02 09:00 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		86		10.0

Comments for Sample AE06986 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 04-22-2002 Time: 09:01:13

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recovery for 7 of the 43 compounds in the LCS Standard were low.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\APR1802\6986.D
 Acq On : 18 Apr 2002 11:41 pm
 Sample : GC/MS SCAN 3/25
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Apr 19 9:00 2002

Vial: 13
 Operator:
 Inst : 209
 Multiplr: 1.00

Quant Results File: GCMS0412.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Last Update : Wed Apr 17 11:48:30 2002
 Response via : Initial Calibration
 DataAcq Meth : GCMS0412

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.21	152	572057	20.00	ug/l	-0.03
10) Naphthalene-d8	15.84	136	2075485	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.15	164	1154037	20.00	ug/l	-0.02
30) Phenanthrene-d10	25.58	188	1689663	20.00	ug/l	-0.03
39) Chrysene-d12	33.64	240	1120948	20.00	ug/l	-0.03
45) Perylene-d12	37.88	264	746032	20.00	ug/l	-0.04

System Monitoring Compounds

28) TCMX	0.00	209	0	0.00	ug/L	
Spiked Amount	10.000		Recovery	=	0.00%	
38) 2,4-DDD	30.66	235	84834	4.31	ug/mL	0.16
Spiked Amount	5.000		Recovery	=	86.20%	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) N-nitroso-dimethyl amine	5.48	74	311	N.D.		
3) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
4) 1,3-Dichlorobenzene	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-Chloropropan	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)methane	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.		
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	22.63	149	686	N.D.		
26) 4-Chlorophenyl-phenylether	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	168	0	N.D.		
32) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		

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

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Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\APR1802\6986.D

Vial: 13

Acq On : 18 Apr 2002 11:41 pm

Operator:

Sample : GC/MS SCAN 3/25

Inst : 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Apr 19 9:00 2002

Quant Results File: GCMS0412.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Wed Apr 17 11:48:30 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0412

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
33) Hexachlorobenzene	0.00	284	0	N.D.	
34) Phenanthrene	25.59	178	464	N.D.	
35) Anthracene	25.59	178	464	N.D.	
36) Di-n-butylphthalate	27.55	149	1557	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	33.64	228	2665	N.D.	
43) Bis(2-ethylhexyl)phthalate	33.86	149	76497	2.71 ug/l	99
44) Chrysene	33.64	228	2666	N.D.	
46) Di-n-Octylphthalate	0.00	149	0	N.D.	
47) Benzo(b)fluoranthene	36.76	252	208	N.D.	
48) Benzo(k)fluoranthene	36.76	252	208	N.D.	
49) Benzo(a)pyrene	37.88	252	3079	N.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

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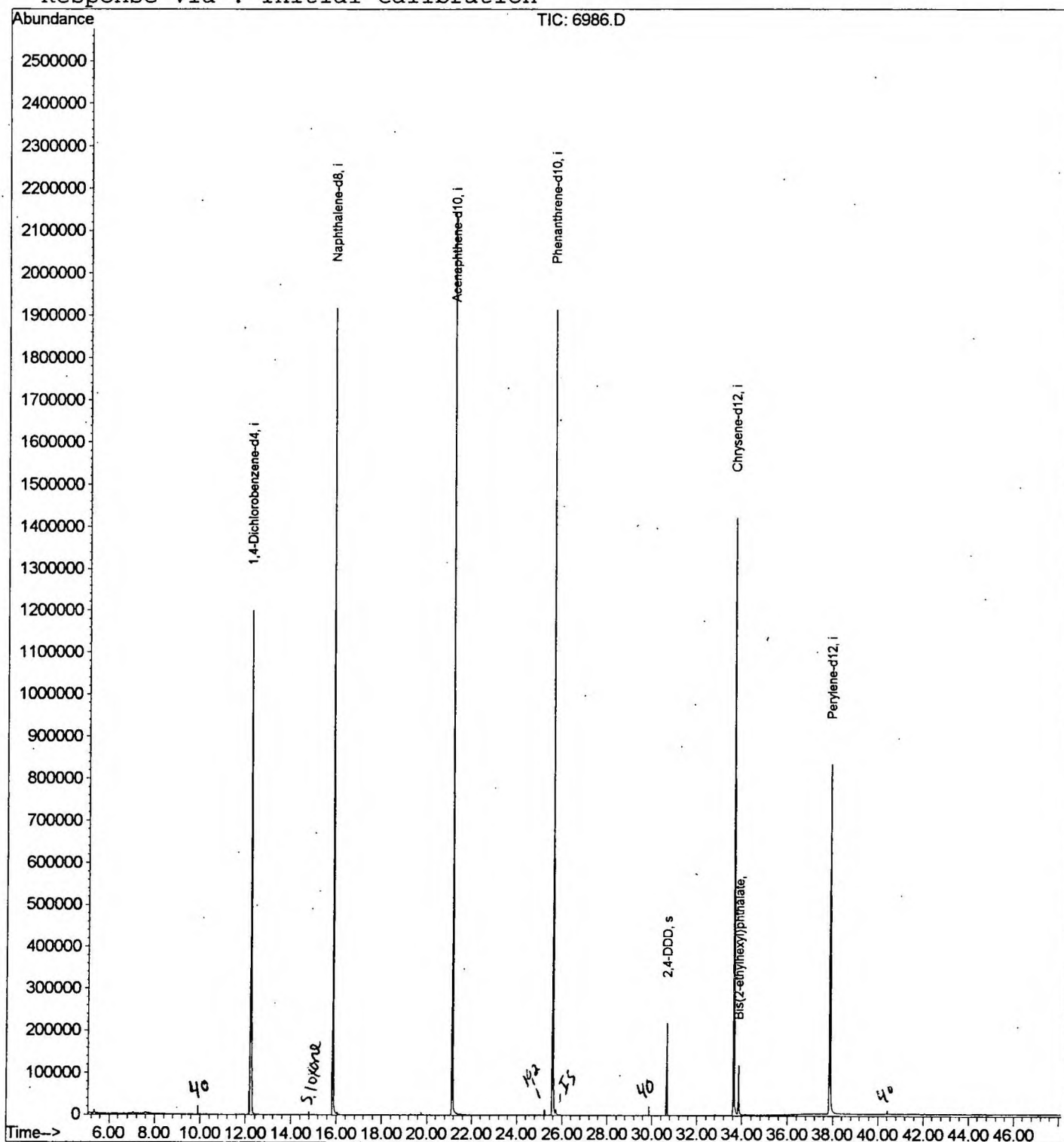
Quantitation Report

Data File : C:\HPCHEM\1\DATA\APR1802\6986.D
 Acq On : 18 Apr 2002 11:41 pm
 Sample : GC/MS SCAN 3/25
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Apr 19 9:00 2002

Vial: 13
 Operator:
 Inst : 209
 Multiplr: 1.00

Quant Results File: GCMS0412.RES

Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Last Update : Wed Apr 17 11:48:30 2002
 Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\APR1802\6986.D
 Acq On : 18 Apr 2002 11:41 pm
 Sample : GC/MS SCAN 3/25
 Misc :
 MS Integration Params: LSCINT.P

Vial: 13
 Operator:
 Inst : 209
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Signal : TIC

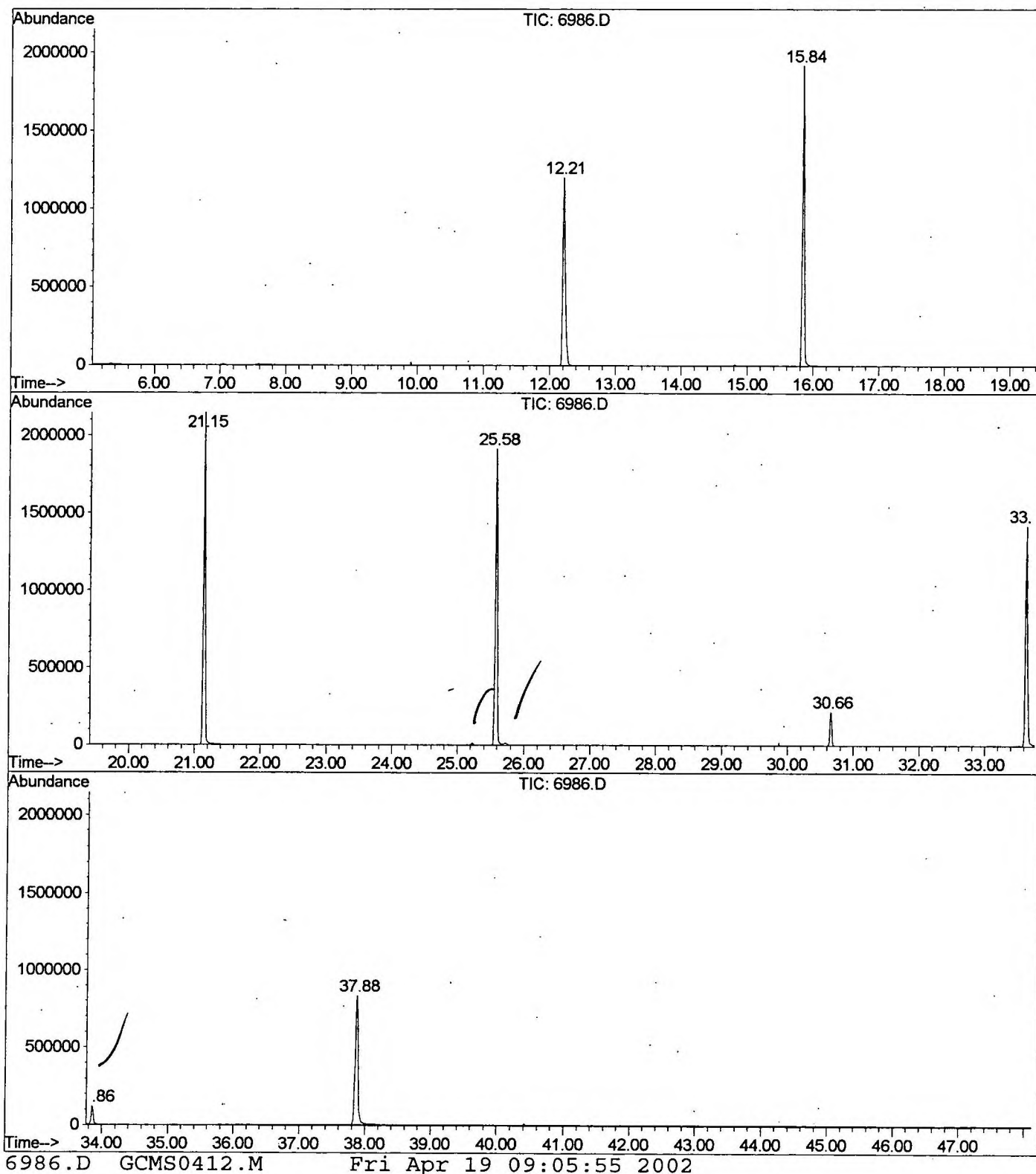
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	12.209	775	780	797	rVB	1198420	3051991	70.31%	13.955%
2	15.844	1170	1176	1194	rBV	1918588	3900214	89.85%	17.833%
3	21.150	1747	1754	1766	rBV	2146041	4340943	100.00%	19.848%
4	25.584	2231	2237	2248	rBV	1914067	4216339	97.13%	19.278%
5	30.661	2785	2790	2797	rBV	218717	414341	9.54%	1.895%
6	33.645	3108	3115	3134	rBV2	1422194	3185949	73.39%	14.567%
7	33.856	3134	3138	3149	rVB	117568	252475	5.82%	1.154%
8	37.877	3566	3576	3594	rBV2	831212	2508428	57.79%	11.469%

Sum of corrected areas: 21870680

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LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\APR1802\6986.D
 Operator :
 Acquired : 18 Apr 2002 11:41 pm using AcqMethod GCMS0412
 Instrument : 209
 Sample Name: GC/MS SCAN 3/25
 Misc Info :
 Vial Number: 13
 Quant File :GCMS0412.RES (RTE Integrator)



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 18 Apr 2002 11:41 pm
Data File: C:\HPCHEM\1\DATA\APR1802\6986.D
Name: GC/MS SCAN 3/25
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0412.M (RTE Integrator)
Title: 209 625/TCLP calibration table
Library Searched: C:\DATABASE\NBS75K.L

TIC	Top	Hit	name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc

6986.D			GCMS0412.M	Fri Apr 19 09:05:55 2002							