

Comments for Sample AD28049 - Analysis \$GCMS

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 02-15-2002 Time: 12:35:32

A scan of the extract prepared for the 508 analysis, from 35-550 amu using a mass spectrometer, does not reveal the presence of non-target compounds. This sample is the Field Blank.

User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/15/2002 Time: 12:35:34

Sample: AD28049
Analysis: \$GCMS GCMS Scan For Unknowns
Units: ug
Started: 1/2/02 15:21 Ended: 1/2/02 15:21 Analyst: MG
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Sum 2,4 DDD	LSPEC	< MDL		5.0

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB0802\28049.D

Vial: 24

Acq On : 9 Feb 2002 7:41 am

Operator: 209 GALOS

Sample : GC/MS SCAN 1/28 FB

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Feb 13 8:49 2002

Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Fri Feb 08 15:29:22 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0208

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	12.35	152	332831	20.00	ug/l	0.00
10) Naphthalene-d8	15.99	136	1271880	20.00	ug/l	0.00
17) Acenaphthene-d10	21.31	164	678039	20.00	ug/l	0.00
29) Phenanthrene-d10	25.76	188	984132	20.00	ug/l	0.00
38) Chrysene-d12	33.83	240	652874	20.00	ug/l	-0.02
44) Perylene-d12	38.11	264	422605	20.00	ug/l	-0.02

System Monitoring Compounds

37) 2,4-DDD	30.83	235	40308	3.29	ug/mL	0.34
Spiked Amount	5.000		Recovery	=	65.80%	

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-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	16.05	128	461	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	0.00	149	0	N.D.
26) 4-Chlorophenyl-phenylether	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.

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

28049.D GCMS0208.M

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Quantitation Report

(QT Reviewed)

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Vial: 24

Acq On : 9 Feb 2002 7:41 am

Operator: 209 GALOS

Sample : GC/MS SCAN 1/28 FB

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Feb 13 8:49 2002

Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Fri Feb 08 15:29:22 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0208

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
30) N-Nitrosodiphenylamine	0.00	168	0	N.D.		
31) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		
32) Hexachlorobenzene	0.00	284	0	N.D.		
33) Phenanthrene	25.76	178	302	N.D.		
34) Anthracene	25.76	178	302	N.D.		
35) Di-n-butylphthalate	27.71	149	16114	N.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	33.82	228	1634	N.D.		
42) Bis(2-ethylhexyl)phthalate	34.03	149	196	N.D.		
43) Chrysene	33.82	228	1634	N.D.		
45) Di-n-Octylphthalate	0.00	149	0	N.D.		
46) Benzo(b)fluoranthene	0.00	252	0	N.D.		
47) Benzo(k)fluoranthene	0.00	252	0	N.D.		
48) Benzo(a)pyrene	38.11	252	1735	N.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

28049.D GCMS0208.M

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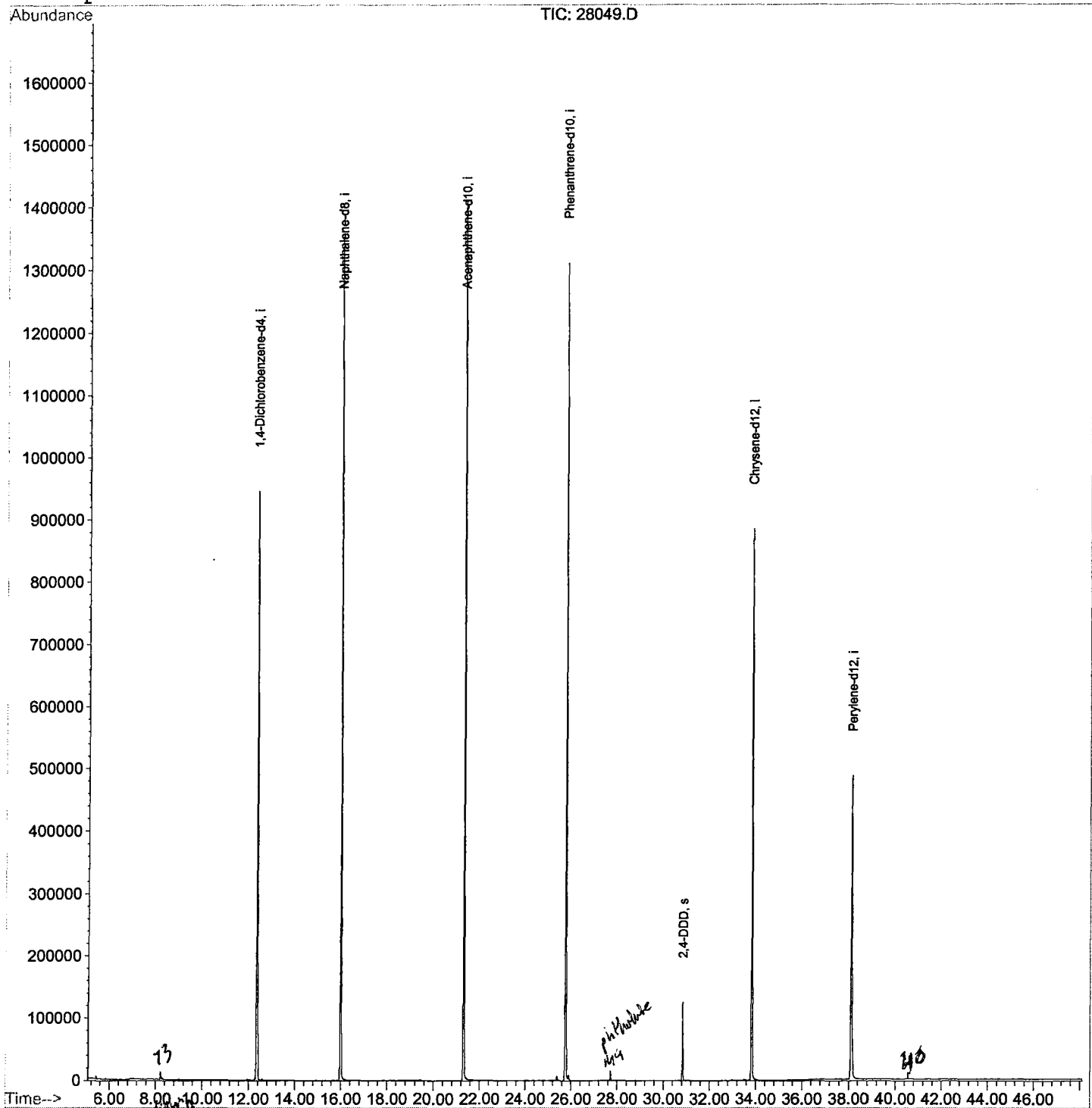
Quantitation Report

Data File : C:\HPCHEM\1\DATA\FEB0802\28049.D
Acq On : 9 Feb 2002 7:41 am
Sample : GC/MS SCAN 1/28 FB
Misc :
MS Integration Params: GOOFY.P
Quant Time: Feb 13 8:49 2002

Vial: 24
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

Quant Results File: GCMS0208.RES

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri Feb 08 15:29:22 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\FEB0802\28049.D

Vial: 24

Acq On : 9 Feb 2002 7:41 am

Operator: 209 GALOS

Sample : GC/MS SCAN 1/28 FB

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)

Title : 209 625/TCLP calibration table

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	peak area	peak % max.	% of total
1	12.346	789	795	810	rVB	944880	2120699	75.44%	15.187%
2	15.990	1186	1192	1208	rBV	1352095	2625909	93.41%	18.805%
3	21.306	1765	1771	1786	rBV	1410598	2811025	100.00%	20.131%
4	25.758	2249	2256	2267	rBV	1310592	2729517	97.10%	19.547%
5	30.835	2804	2809	2814	rVB	124967	239504	8.52%	1.715%
6	33.828	3128	3135	3148	rBV	884132	1984041	70.58%	14.208%
7	38.115	3593	3602	3616	rBV2	486247	1453248	51.70%	10.407%

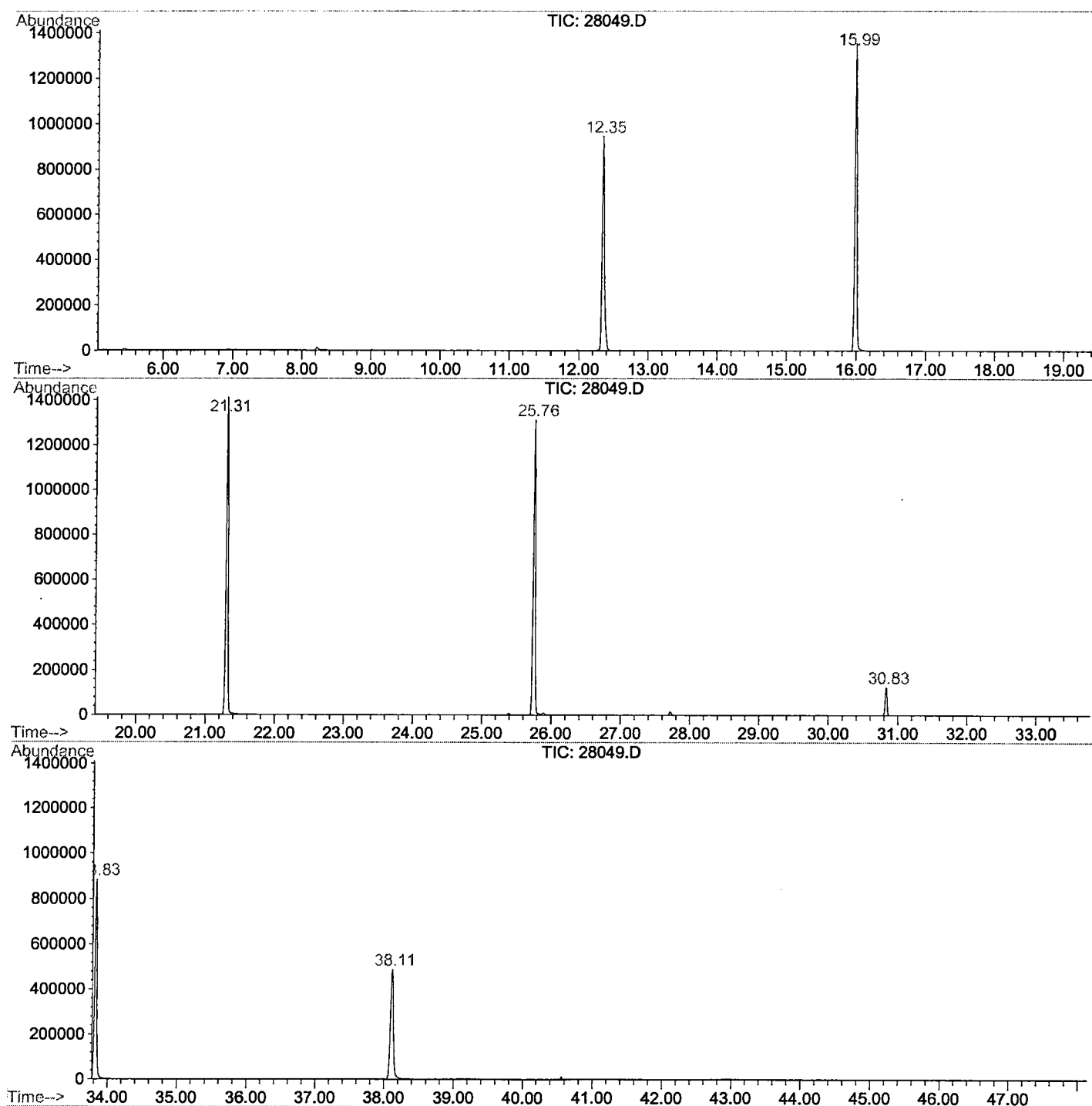
Sum of corrected areas: 13963943

28049.D GCMS0208.M

Thu Feb 14 15:02:12 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\FEB0802\28049.D
Operator : 209 GALOS
Acquired : 9 Feb 2002 7:41 am using AcqMethod GCMS0208
Instrument : GC/MS 209
Sample Name: GC/MS SCAN 1/28 FB
Misc Info :
Vial Number: 24
Quant File :GCMS0208.RES (RTE Integrator)



28049.D GCMS0208.M

Thu Feb 14 15:02:18 2002

Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 9 Feb 2002 7:41 am
 Data File: C:\HPCHEM\1\DATA\FEB0802\28049.D
 Name: GC/MS SCAN 1/28 FB
 Misc:
 Method: C:\HPCHEM\1\METHODS\GCMS0208.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

28049.D GCMS0208.M	Thu Feb 14	15:02:22	2002					