

User: Fakhouri, Morgan
 Westchester Dept. of Labs & Research
 Date: 01/29/2002 Time: 11:34:09

Sample: AD28166
 Analysis: \$525 Organic Chemicals - 525.2
 Units: ug
 Started: 1/29/02 11:33 Ended: 1/29/02 11:33 Analyst: MF
 Result source: manual entry
 Page: 1

original data

	Component Name	Viol	Result	Qualifier	MDL
1	Hexachlorocyclopentadiene		< MDL		0.10
2	Hexachlorobenzene		< MDL		0.10
3	Simazine		< MDL		0.40
4	Atrazine		< MDL		0.20
5	Alachlor		< MDL		0.20
6	bis(2-Ethylhexyl)adipate		< MDL		0.60
7	bis(2-Ethylhexyl)phthalate		< MDL		0.60
8	Benzo(a)pyrene		< MDL		0.20
9	EPTC		< MDL		1.0
10	2,6-Dinitrotoluene		< MDL		2.0
11	2,4-Dinitrotoluene		< MDL		2.0
12	Molinate		< MDL		0.90
13	Terbacil		< MDL		2.0
14	Acetochlor		< MDL		2.0
15	Metribuzin		< MDL		0.15
16	Metolachlor		< MDL		0.75
17	Butachlor		< MDL		0.40
18	4,4-DDE		< MDL		0.80
19	Surr_1,3-Dimethyl-2-nitrobenzen		4.31		1.0
20	Surr_Triphenyl phosphate		6.21		1.0
21	Surr_Perylene-d12		4.06		1.0

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\121001B\28166.D

Acq On : 11 Dec 2001 8:42 am

Sample : sample

Misc :

MS Integration Params: BENZO.P

Quant Time: Dec 11 14:11:10 2001

Vial: 23

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

Quant Results File: 120501.RES

Original data

Quant Method : C:\MSDCHEM\1\5973N\120501.M (RTE Integrator)

Title : Quantitation For Method 525.2

Last Update : Thu Dec 06 10:31:04 2001

Response via : Initial Calibration

DataAcq Meth : 120501

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Acenaphthene-d10	18.32	164	6649733	5.00	ug/L	0.00
9) Phenanthrene D-10	22.66	188	10254163	5.00	ug/L	0.02
20) Chrysene D-12	30.43	240	6427829	5.00	ug/L	0.00

System Monitoring Compounds

2) 1,3-Dimethyl-2-nitrobenzen	13.24	134	1519893	4.31	ug/L	0.02
Spiked Amount	5.000		Recovery	=	86.20%	
19) Triphenyl Phosphate surr	29.70	326	2207223	6.21	ug/L	0.00
Spiked Amount	5.000		Recovery	=	124.20%	
21) Perylene D-12 surr	34.45	264	3659403	4.06	ug/L	0.00
Spiked Amount	5.000		Recovery	=	81.20%	

Target Compounds

					Qvalue
3) Hexachlorocyclopentadiene	0.00	237	0	N.D.	
4) Hexachlorobenzene	21.52	284	7304	0.01 ug/L #	75
5) EPTC	0.00	128	0	N.D.	
6) 2,6-Dinitrotoluene	0.00	165	0	N.D. d	
7) 2,4-Dinitrotoluene	0.00	165	0	N.D. d	
8) Molinate	0.00	126	0	N.D. d	
10) Simazine	22.10	201	6771	0.02 ug/L #	18
11) Atrazine	0.00	200	0	N.D.	
12) Alachlor	0.00	160	0	N.D. d	
13) Terbacil	0.00	161	0	N.D. d	
14) Acetochlor	24.23	146	10117	0.04 ug/L #	79
15) Metribuzin	0.00	198	0	N.D.	
16) Metolachlor	25.38	162	5826	0.01 ug/L	67
17) Butachlor	26.58	160	6473	0.02 ug/L	92
18) 4,4'-DDE	0.00	246	0	N.D. d	
22) Bis (2-Ethylhexyl) adipate	29.62	129	42690	0.15 ug/L #	21
23) Bis (2-Ethylhexyl) phthala	31.07	149	429709	0.43 ug/L	90
24) Benzo (a) pyrene	0.00	252	0	N.D. d	

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

28166.D 120501.M Thu Jan 10 12:14:41 2002

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\121001B\28166.D

Vial: 23

Acq On : 11 Dec 2001 8:42 am

Operator: McLean

Sample : sample

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: BENZO.P

Quant Time: Jan 10 12:14 2002

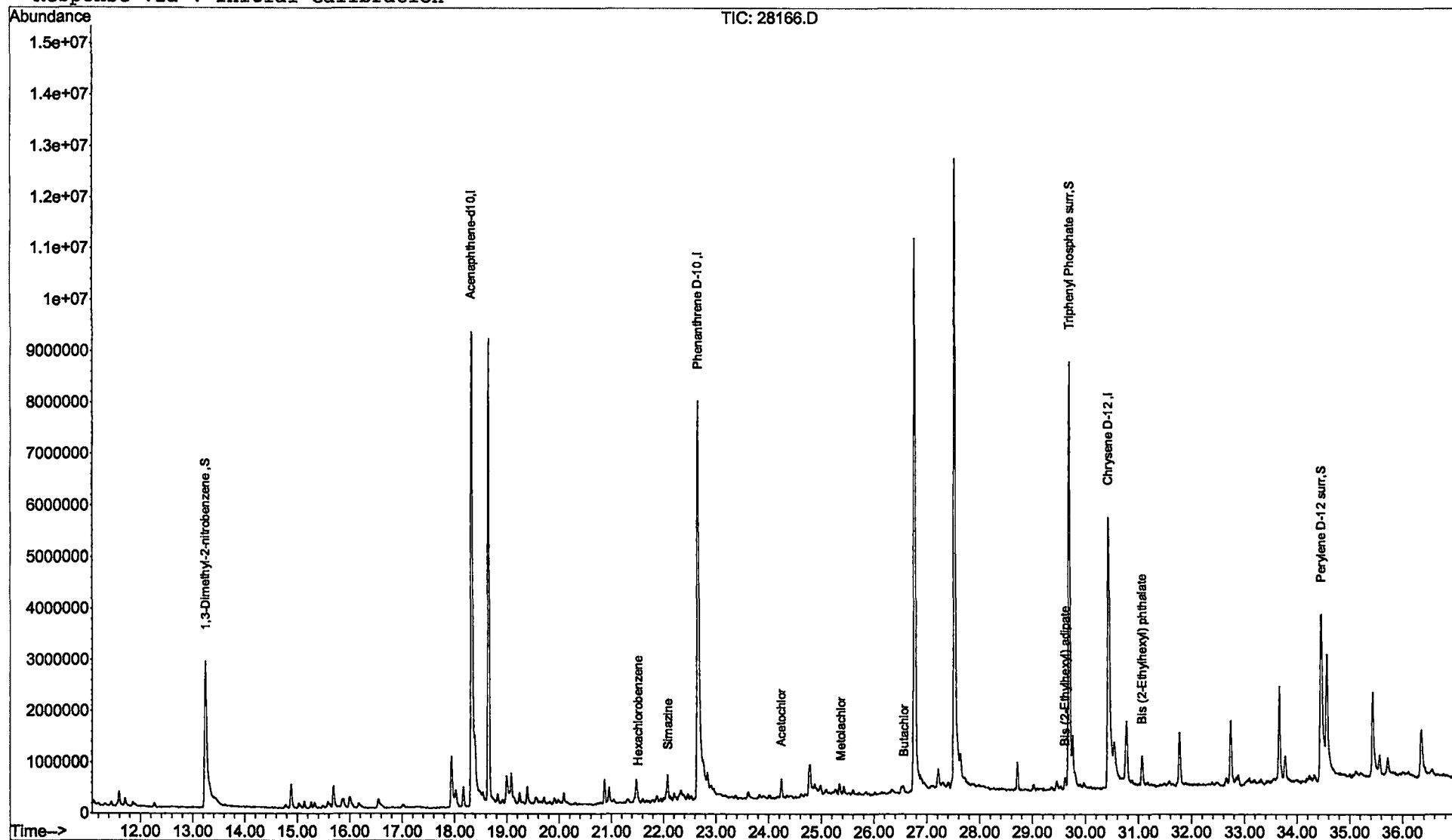
Quant Results File: 120501.RES

Method : C:\MSDCHEM\1\5973N\120501.M (RTE Integrator)

Title : Quantitation For Method 525.2

Last Update : Thu Dec 06 10:31:04 2001

Response via : Initial Calibration



Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\012502A\28166RR.D

Acq On : 25 Jan 2002 7:10 pm

Sample : REINJ

Misc :

MS Integration Params: rteint.p

Quant Time: Jan 28 12:40:02 2002

Vial: 7

Operator: McLean

Inst : Instrumen

Multiplr: 1.00

Quant Results File: 012402B.RES

revised data

Quant Method : C:\MSDCHEM\1\5973N\012402B.M (RTE Integrator)

Title : Quantitation For Method 525.2

Last Update : Fri Jan 25 11:19:14 2002

Response via : Initial Calibration

DataAcq Meth : 012402DF

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Acenaphthene-d10	17.96	164	14263485	5.00	ug/L	0.00
9) Phenanthrene D-10	22.27	188	18670009	5.00	ug/L	0.00
20) Chrysene D-12	30.03	240	13361418	5.00	ug/L	0.00

System Monitoring Compounds

2) 1,3-Dimethyl-2-nitrobenzen	12.89	134	3709701	5.07	ug/L	0.00
Spiked Amount	5.000		Recovery	=	101.40%	
19) Triphenyl Phosphate	29.33	326	3370287	5.99	ug/L	0.00
Spiked Amount	5.000		Recovery	=	119.80%	
21) Perylene D-12	34.02	264	7206213	3.80	ug/L	0.00
Spiked Amount	5.000		Recovery	=	76.00%	

Target Compounds

					Qvalue
3) Hexachlorocyclopentadiene	0.00	237	0	N.D. d	
4) Hexachlorobenzene	21.14	284	16149	0.02 ug/L	83
5) EPTC	0.00	128	0	N.D.	
6) 2,6-Dinitrotoluene	0.00	165	0	N.D. d	
7) 2,4-Dinitrotoluene	18.70	165	72739	0.98 ug/L	27
8) Molinate	18.90	126	5657	Below Cal #	1
10) Simazine	0.00	201	0	N.D. d	
11) Atrazine	0.00	200	0	N.D.	
12) Alachlor	0.00	160	0	N.D. d	
13) Terbacil	0.00	161	0	N.D. d	
14) Acetochlor	0.00	146	0	N.D. d	
15) Metribuzin	0.00	198	0	N.D.	
16) Metolachlor	24.59	162	6715	0.22 ug/L	81
17) Butachlor	0.00	160	0	N.D. d	
18) 4,4'-DDE	27.05	246	9571	0.04 ug/L	96
22) Bis (2-Ethylhexyl) adipate	29.27	129	60671	0.53 ug/L #	1
23) Bis (2-Ethylhexyl) phthala	30.69	149	612843	0.60 ug/L	96
24) Benzo (a) pyrene	0.00	252	0	N.D. d	

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

28166RR.D 012402B.M

Mon Jan 28 12:43:58 2002

Page 1

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\012502A\28166RR.D

Vial: 7

Acq On : 25 Jan 2002 7:10 pm

Operator: McLean

Sample : REINJ

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: rteint.p

Quant Time: Jan 28 12:43 2002

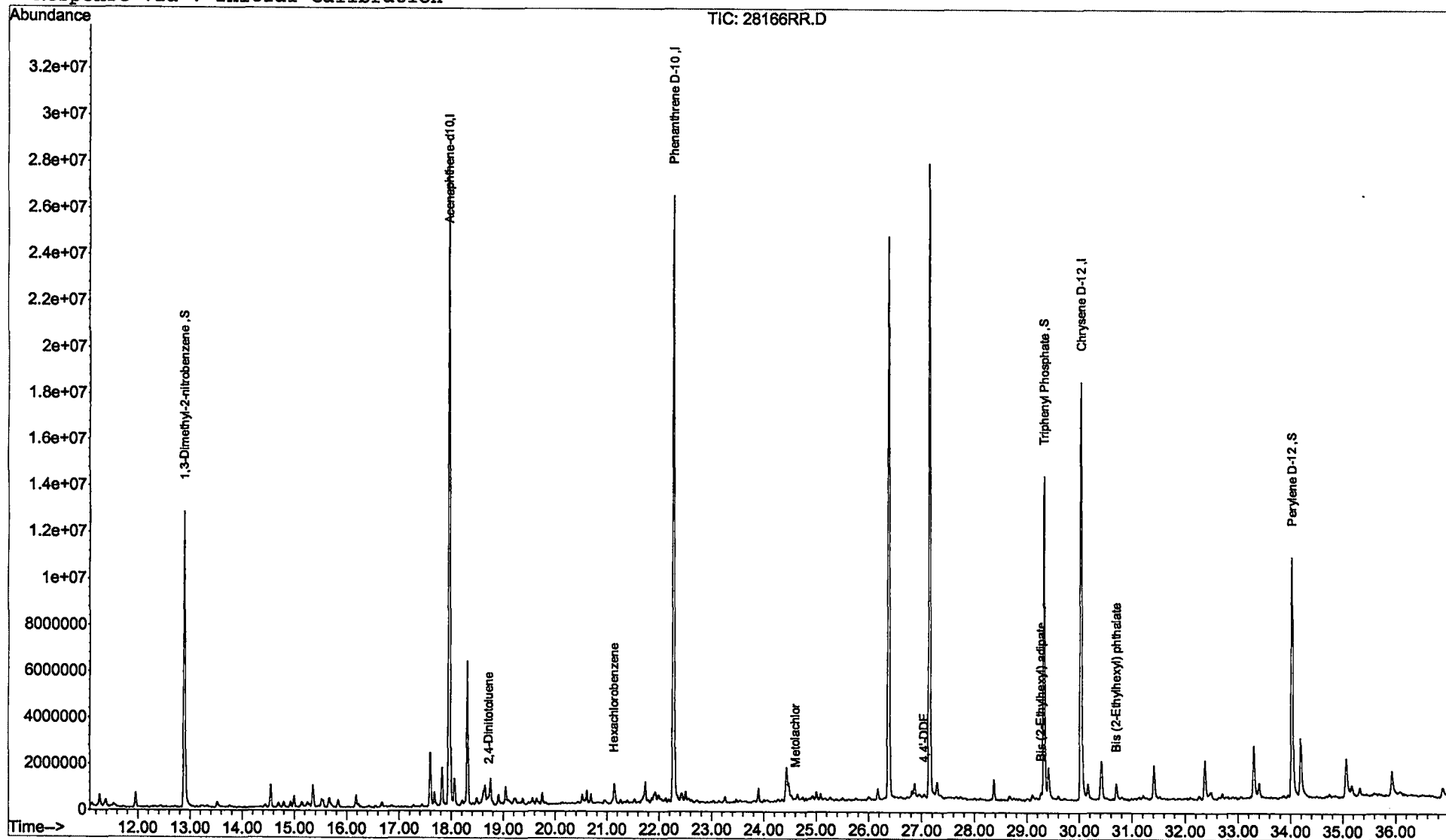
Quant Results File: 012402B.RES

Method : C:\MSDCHEM\1\5973N\012402B.M (RTE Integrator)

Title : Quantitation For Method 525.2

Last Update : Fri Jan 25 11:19:14 2002

Response via : Initial Calibration



Comments for Sample AD28166 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 01-30-2002 Time: 15:09:06

The recoveries for 6 of the 18 compounds in the LCS Standard were outside of their acceptance ranges. The recoveries for 4 of the 18 compounds in the Spiked Sample were outside of their acceptance ranges.