re-extracted
(original data)

User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 01/29/2002 Time: 10:12:57

Sample: AD28043
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 1/29/02 10:10 Ended: 1/29/02 10:10 Analyst: MF
Result source: manual entry
Page: 1

	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		5.14		1.0
20	Surr_Triphenyl phosphate		6.12		1.0
21	Surr_Perylene-d12		4.85		1.0

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\012502A\28043RR.D Vial: 14
Acq On : 26 Jan 2002 12:25 am Operator: McLean
Sample : REINJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jan 28 11:56:54 2002 Quant Results File: 012402B.RES

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

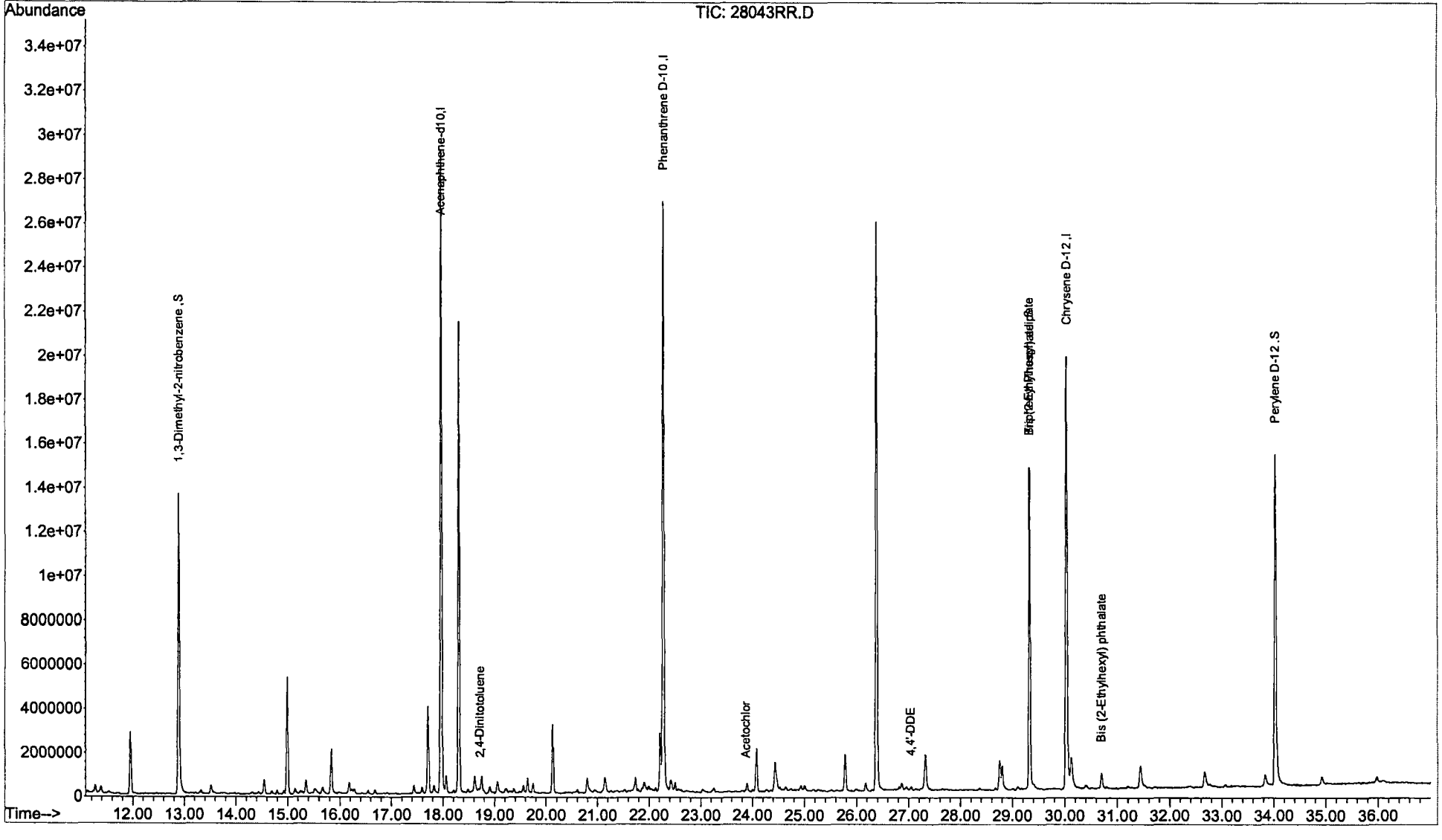
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Acenaphthene-d10	17.96	164	14887214	5.00	ug/L	0.00
9) Phenanthrene D-10	22.27	188	19442542	5.00	ug/L	0.00
20) Chrysene D-12	30.03	240	14693109	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	12.89	134	3920217	5.14	ug/L	0.00
Spiked Amount 5.000			Recovery = 102.80%			
19) Triphenyl Phosphate	29.33	326	3585371	6.12	ug/L	0.00
Spiked Amount 5.000			Recovery = 122.40%			
21) Perylene D-12	34.02	264	10128960	4.85	ug/L	0.00
Spiked Amount 5.000			Recovery = 97.00%			
Target Compounds						Qvalue
3) Hexachlorocyclopentadiene	0.00	237	0	N.D.	d	
4) Hexachlorobenzene	0.00	284	0	N.D.	d	
5) EPTC	15.82	128	7112	N.D.		
6) 2,6-Dinitrotoluene	0.00	165	0	N.D.	d	
7) 2,4-Dinitotoluene	18.70	165	6436	0.91	ug/L	45
8) Molinate	18.90	126	15131	Below Cal	#	1
10) Simazine	0.00	201	0	N.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	23.86	146	6427	0.43	ug/L #	41
15) Metribuzin	0.00	198	0	N.D.		
16) Metolachlor	0.00	162	0	N.D.		
17) Butachlor	0.00	160	0	N.D.	d	
18) 4,4'-DDE	27.03	246	15975	0.05	ug/L	96
22) Bis (2-Ethylhexyl) adipate	29.32	129	66103	0.53	ug/L	25
23) Bis (2-Ethylhexyl) phthala	30.69	149	610548	0.58	ug/L	99
24) Benzo (a) pyrene	0.00	252	0	N.D.	d	

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

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\012502A\28043RR.D Vial: 14
Acq On : 26 Jan 2002 12:25 am Operator: McLean
Sample : REINJ Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Jan 28 11:59 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



Quantitation Report (Not Reviewed)

re-extracted reinjection Data

Data File : C:\MSDCHEM\1\DATA\121001B\28043.D Vial: 11
Acq On : 10 Dec 2001 11:54 pm Operator: McLean
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 11 00:31:39 2001 Quant Results File: 120501.RES

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	18.44	164	141697	5.00	ug/L	0.13
9) Phenanthrene D-10	22.80	188	24122	5.00	ug/L	0.16
20) Chrysene D-12	30.56	240	81288	5.00	ug/L	0.14
System Monitoring Compounds						
2) 1,3-Dimethyl-2-nitrobenzen	13.36	134	5022	0.67	ug/L	0.14
Spiked Amount 5.000			Recovery = 13.40%			
19) Triphenyl Phosphate surr	29.75	326	62350	74.56	ug/L	0.06
Spiked Amount 5.000			Recovery = 1491.20%			
21) Perylene D-12 surr	34.54	264	58078	5.09	ug/L	0.10
Spiked Amount 5.000			Recovery = 101.80%			
Target Compounds						Qvalue
3) Hexachlorocyclopentadiene	0.00	237	0	N.D.		
4) Hexachlorobenzene	0.00	284	0	N.D.		
5) EPTC	0.00	128	0	N.D.		
6) 2,6-Dinitrotoluene	18.44	165	13959	2.70	ug/L #	30
7) 2,4-Dinitotoluene	19.06	165	1221	0.47	ug/L #	21
8) Molinate	0.00	126	0	N.D.		
10) Simazine	0.00	201	0	N.D.		
11) Atrazine	0.00	200	0	N.D.		
12) Alachlor	0.00	160	0	N.D.		
13) Terbacil	0.00	161	0	N.D.		
14) Acetochlor	0.00	146	0	N.D.		
15) Metribuzin	0.00	198	0	N.D.		
16) Metolachlor	0.00	162	0	N.D.		
17) Butachlor	0.00	160	0	N.D.		
18) 4,4'-DDE	0.00	246	0	N.D.		
22) Bis (2-Ethylhexyl) adipate	0.00	129	0	N.D		
23) Bis (2-Ethylhexyl) phthala	31.08	149	17174	0		78
24) Benzo (a) pyrene	0.00	252	0			

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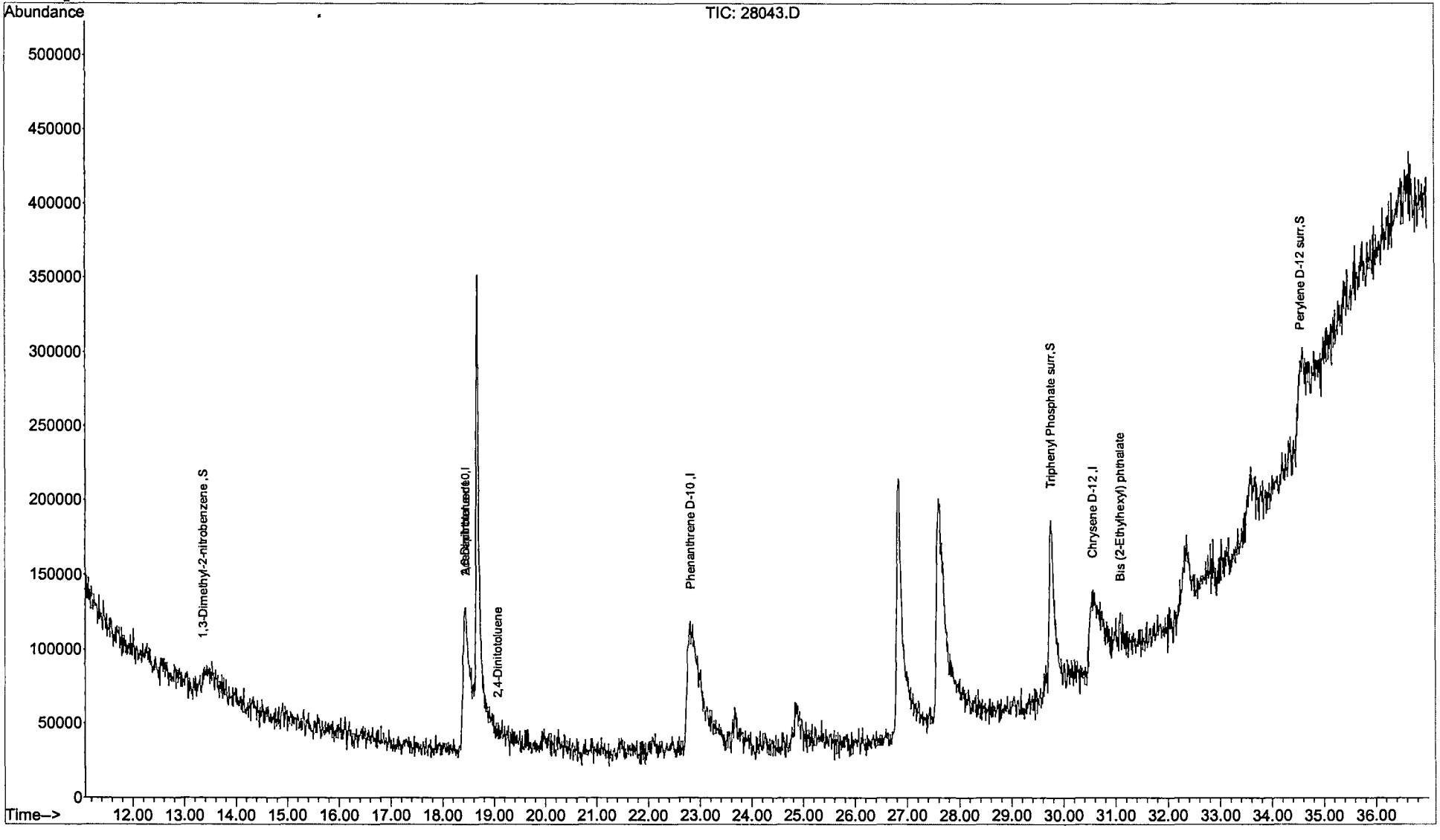
the old needed to be re-extracted

(#) = qualifier out of range (m) = m
28043.D 012402B.M Mon Jan 28 1

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\DATA\121001B\28043.D Vial: 11
Acq On : 10 Dec 2001 11:54 pm Operator: McLean
Sample : sample Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: BENZO.P
Quant Time: Dec 11 0:31 2001 Quant Results File: 120501.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 AD28043 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 01-30-2002 Time: 15:26:09

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.