

User: Bohlk, Ted
 Westchester Dept. of Labs & Research
 Date: 12/13/2001 Time: 18:33:41

Sample: AD26079
 Analysis: \$525 Organic Chemicals - 525.2
 Units: ug
 Started: 11/1/01 18:32 Ended: 11/17/01 18:32 Analyst: TB
 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		0.88		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-nitrobenze		4.7		1.0
20	Surr_Triphenyl phosphate		5.7		1.0
21	Surr_Perylene d12	LSPEC	3.3		1.0

Comments for Sample AD26079 - Analysis \$525

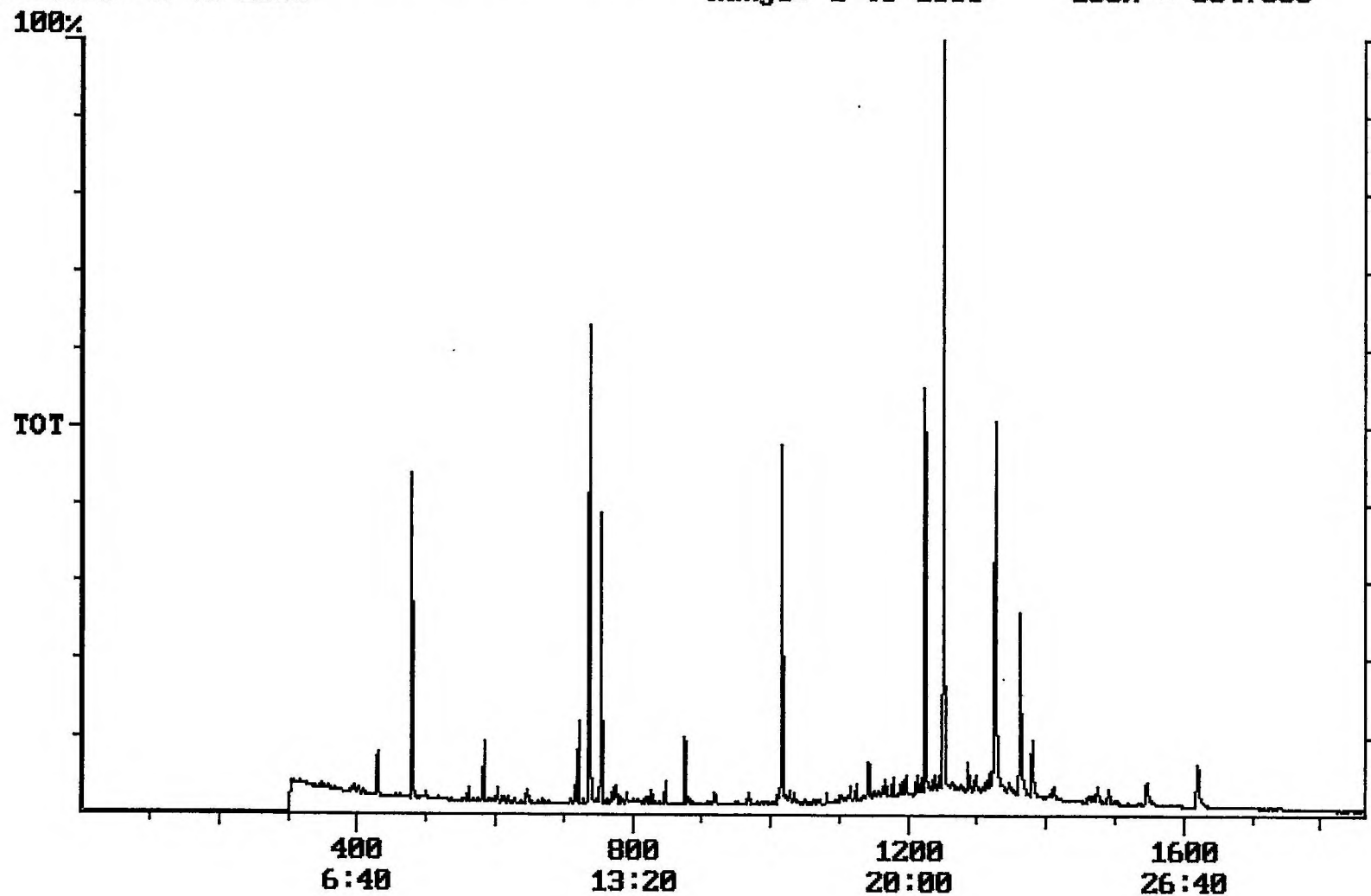
Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 12-15-2001 Time: 11:43:43

The recovery for 1 of the 3 Surrogate Standards in this sample is below its acceptance range. Recoveries for 3 of the 18 compounds in the MDL and LCS Standards are outside of their acceptance ranges. Recoveries for 4 of the 18 compounds in the Spiked Sample are outside of their acceptance ranges. The level for bis(2-Ethylhexyl)phthalate in the Laboratory Blank is 1.2 ug/L.

Chromatogram Plot File: E:\260790 Date: Nov-18-1991 09:30:33
Comment: BOHLK; METHOD 525; INST 204; 11/17/01; SAMPLE 26079
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 8347000



Integration Report

Quanfile: 260790

Quan Entries: 15

Comment: BOHLK; METHOD 525; INST 204; 11/17/01; SAMPLE 26079

Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:18	738	Auto	1,930,783	1,222,839
2	phenanthrene d-10 (IS)	16:56	1016	Auto	3,141,838	1,337,776
3	chrysene d-12(IS)	22:41	1361	Auto	1,580,393	570,674
4	p-terphenyl d-14(IS)	20:51	1251	Auto	3,756,994	2,498,383
5	acenaphthene d-10(SURR)	12:18	738	Auto	1,930,783	1,222,839
6	phenanthrene d-10(SURR)	16:56	1016	Auto	3,141,838	1,337,776
7	chrysene d-12 (SURR)	22:41	1361	Auto	1,580,393	570,674
8	1,3-dimethyl-2-nitrobenzene	7:59	479	Auto	1,139,442	528,658
9	triphenyl phosphate (S)	22:06	1326	Auto	1,298,016	588,780
10	perylene d-12 (SURR)	27:00	1620	Auto	629,993	104,615
11	hexachlorobenzene	15:22	922	Auto	1,114	1,114
12	simazine	16:08	968	Auto	5,694	1,687
13	bis(2-ethylhexyl)adipate	22:03	1323	Auto	52,882	6,232
14	bis(2-ethylhexyl)phthalate	22:59	1379	Auto	536,725	165,838
15	2,6-dinitrotoluene	11:58	718	Auto	48,642	25,715

Quantitation Report

Quanfile: 260790

Quan Entries: 15

Comment: BOHLK; METHOD 525; INST 204; 11/17/01; SAMPLE 26079

Sorted via: Entry Number ↑

(S) = Standard

Cal	Name of Compound	S	Scan#	R Time	Me	Calc Amt(A)	Units
1	acenaphthene d-10(IS)	I	738	12:18	BV	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1016	16:56	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1361	22:41	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1251	20:51	BB	5.000	UG/L
5	acenaphthene d-10(SURR)	A	738	12:18	BV	3.382	UG/L
6	phenanthrene d-10(SURR)	A	1016	16:56	BV	3.589	UG/L
7	chrysene d-12 (SURR)	A	1361	22:41	BB	2.706	UG/L
8	1,3-dimethyl-2-nitrobe	A	479	7:59	BV	4.738	UG/L
9	triphenyl phosphate (S	A	1326	22:06	BV	5.729	UG/L
10	perylene d-12 (SURR)	A	1620	27:00	BB	3.303	UG/L
12	hexachlorobenzene	A	922	15:22	BB	0.005	UG/L
13	simazine	A	968	16:08	BB	0.064	UG/L
16	bis(2-ethylhexyl)adipa	A	1323	22:03	MM	0.278	UG/L
17	bis(2-ethylhexyl)phtha	A	1379	22:59	BB	0.884	UG/L
20	2,6-dinitrotoluene	A	718	11:58	BB	0.388	UG/L