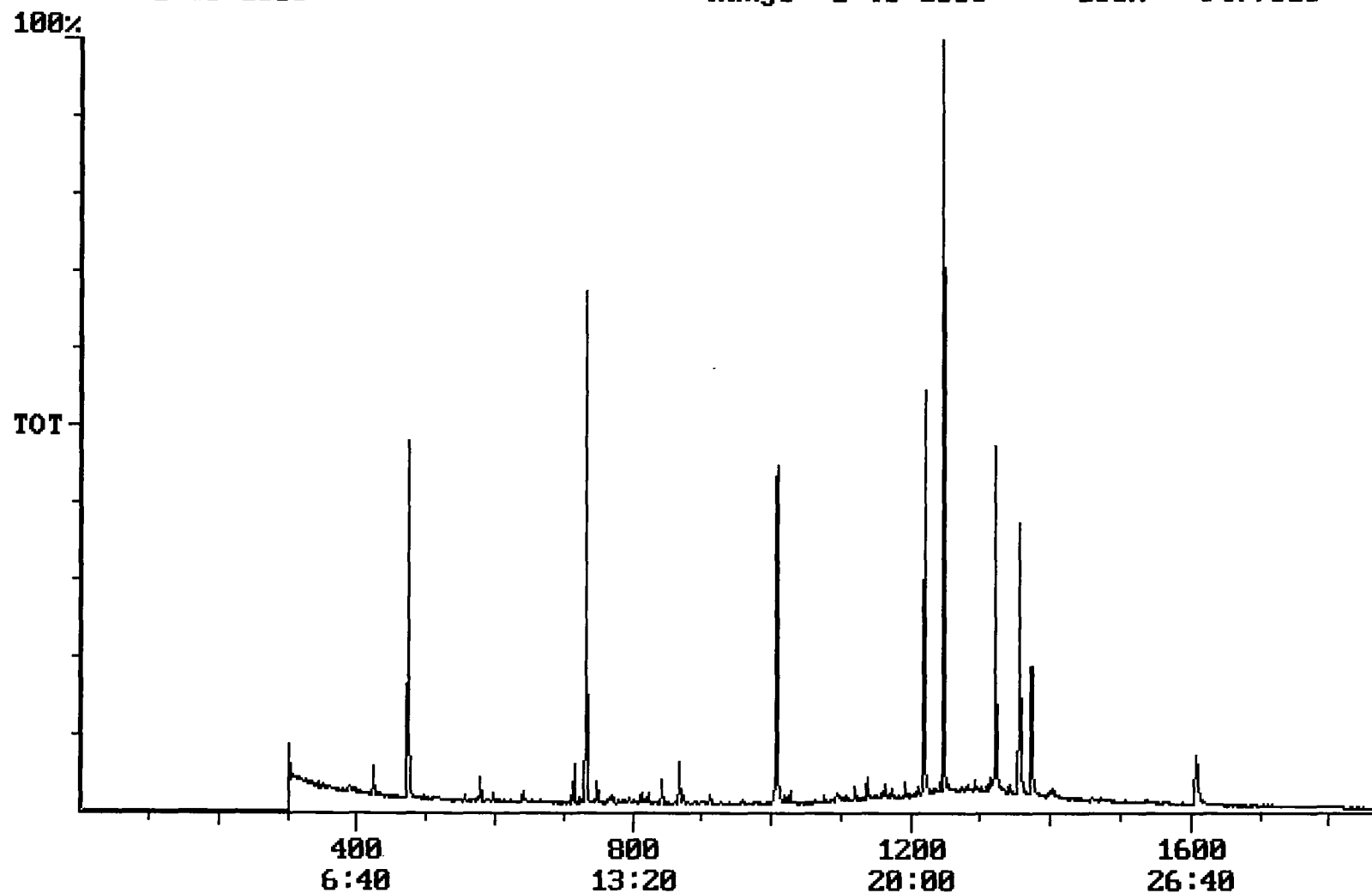


User: Bohik, Ted
 Westchester Dept. of Labs & Research
 Date: 12/19/2001 Time: 19:05:12

Sample: AD26388
 Analysis: \$525 Organic Chemicals - 525.2
 Units: ug
 Started: 11/6/01 19:03 Ended: 11/23/01 19:03 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		1.9		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		4.8		1.0
20	Surr_Triphenyl phosphate		5.4		1.0
21	Surr_Perylene-d12	LSPEC	4.3		1.0

Chromatogram Plot File: E:\26388G Date: Nov-24-1991 05:30:39
Comment: BOHLK; METHOD 525; INST 204; 11/23/01; SAMPLE 26388
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 9477615



Integration Report Quanfile: 26388G Quan Entries: 15
 Comment: BOHLK; METHOD 525; INST 204; 11/23/01; SAMPLE 26388
 Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10 (IS)	12:12	732	Auto	2,397,127	1,573,831
2	phenanthrene d-10 (IS)	16:47	1007	Auto	3,716,245	1,353,566
3	chrysene d-12 (IS)	22:35	1355	Auto	2,155,009	1,021,481
4	p-terphenyl d-14 (IS)	20:46	1246	Auto	4,386,546	2,946,787
5	acenaphthene d-10 (SURR)	12:12	732	Auto	2,397,127	1,573,831
6	phenanthrene d-10 (SURR)	16:47	1007	Auto	3,716,245	1,353,566
7	chrysene d-12 (SURR)	22:35	1355	Auto	2,155,009	1,021,481
8	1,3-dimethyl-2-nitrobenzene	7:54	474	Auto	1,391,552	672,926
9	triphenyl phosphate (S)	22:01	1321	Auto	1,292,814	566,605
10	perylene d-12 (SURR)	26:47	1607	Auto	831,373	171,142
11	hexachlorocyclopentadiene	10:04	604	Auto	6,166	3,479
12	bis(2-ethylhexyl)adipate	21:54	1314	Auto	50,811	11,147
13	bis(2-ethylhexyl)phthalate	22:52	1372	Auto	1,432,603	460,702
14	2,4-dinitrotoluene	13:00	780	Auto	6,110	1,685
15	butachlor	20:19	1219	Auto	1,739	1,739

reversed IS/ SURR's

Quantitation Report

Quanfile: 26388G

Quan Entries: 15

Comment: BOHLK; METHOD 525; INST 204; 11/23/01; SAMPLE 26388

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	732	12:12	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1007	16:47	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1355	22:35	BV	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1246	20:46	BB	5.000	UG/L
5	acenaphthene d-10(SURR)	A	732	12:12	BB	3.604	UG/L
6	phenanthrene d-10(SURR)	A	1007	16:47	BV	3.889	UG/L
7	chrysene d-12 (SURR)	A	1355	22:35	BV	2.935	UG/L
8	1,3-dimethyl-2-nitrobe	A	474	7:54	BV	4.838	UG/L
9	triphenyl phosphate (S	A	1321	22:01	BB	5.411	UG/L
10	perylene d-12 (SURR)	A	1607	26:47	BB	3.342	UG/L
11	hexachlorocyclopentadi	A	604	10:04	BB	0.059	UG/L
16	bis(2-ethylhexyl)adipa	A	1314	21:54	MM	0.105	UG/L
17	bis(2-ethylhexyl)phtha	A	1372	22:52	BV	1.910	UG/L
21	2,4-dinitrotoluene	A	780	13:00	MM	0.039	UG/L
27	butachlor	A	1219	20:19	BB	0.007	UG/L

Comments for Sample AD26388 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 12-21-2001 Time: 16:32:41

The breakdown for Endrin in the Breakdown Standard was 56%. Recoveries for 4 of the 18 compounds in the MDL Standard were outside of their acceptance ranges. Recoveries for 3 of the 18 compounds in the LCS Standard were outside of their acceptance ranges. Recoveries for 3 of the 18 compounds in the Spiked Sample were outside of their acceptance ranges. The level for bis(2-Ethylhexyl)phthalate was 0.88 ug/L in the Laboratory Blank. The recovery for 1 of the 3 Surrogate Standards in this sample was below its acceptance range.