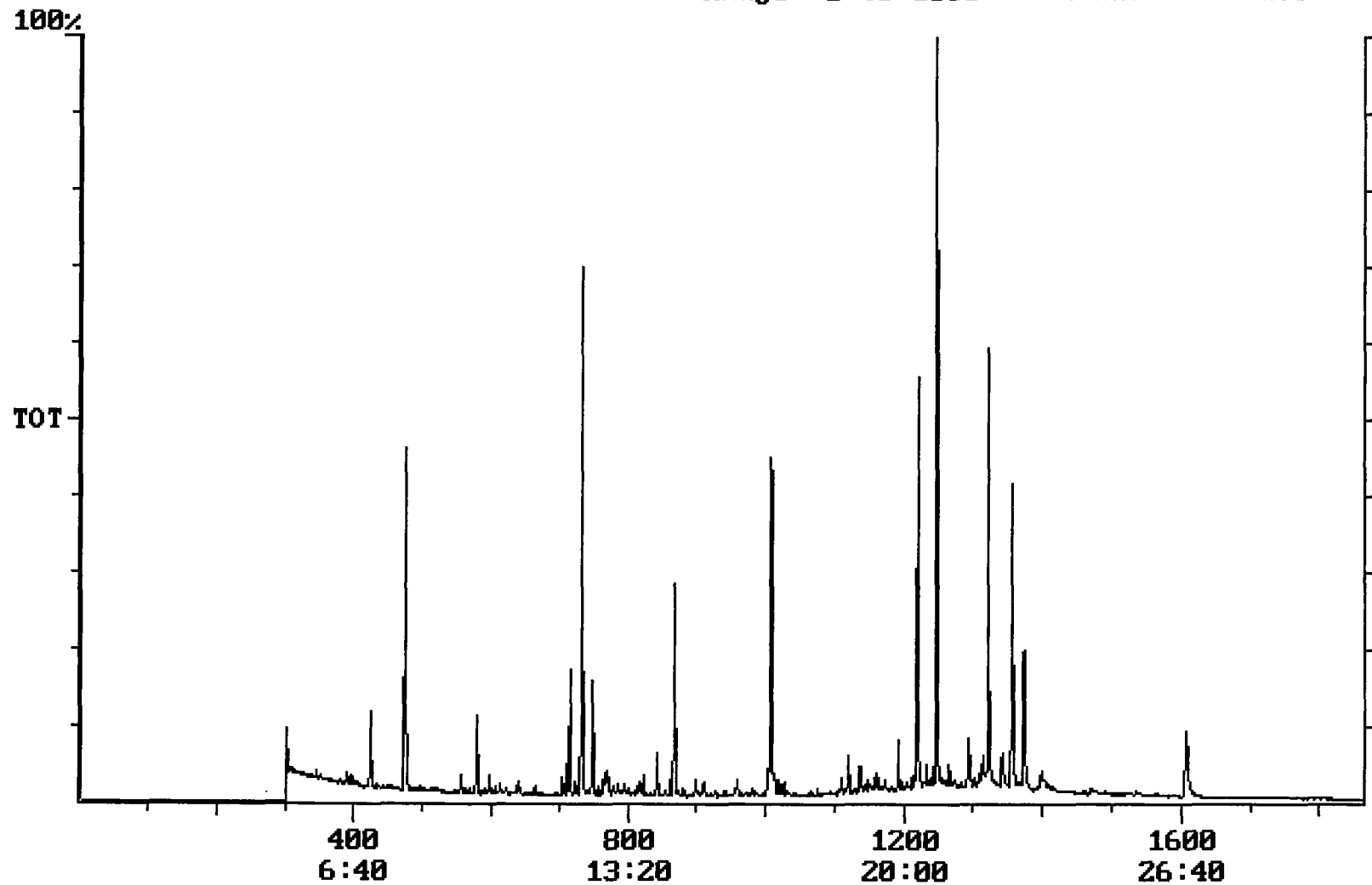


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/20/2001 Time: 09:31:49

Sample: AD26390
Analysis: \$525 Organic Chemicals - 525.2
Units: ug
Started: 11/6/01 09:30 Ended: 11/23/01 09:30 Analyst: TB
Result source: manual entry
Page: 1

	Component Name	Viol	Result	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		1.8	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		4.7	1.0
20	Surr_Triphenyl phosphate		5.5	1.0
21	Surr_Perylene-d12		3.6	1.0

Chromatogram Plot File: E:\26390I Date: Nov-24-1991 12:05:49
Comment: BOHLK; METHOD 525; INST 204; 11/23/01; SAMPLE 26390
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 8426373



Integration Report

Quanfile: 263901

Quan Entries: 17

Comment: BOHLK; METHOD 525; INST 204; 11/23/01; SAMPLE 26390

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	2,211,264	1,416,964
2	phenanthrene d-10 (IS)	16:47	1007	Auto	3,425,252	1,270,365
3	chrysene d-12(IS)	22:35	1355	Auto	2,163,090	1,012,214
4	p-terphenyl d-14(IS)	20:46	1246	Auto	3,825,160	2,594,032
5	acenaphthene d-10(SURR)	12:12	732	Auto	2,211,264	1,416,964
6	phenanthrene d-10(SURR)	16:47	1007	Auto	3,425,252	1,270,365
7	chrysene d-12 (SURR)	22:35	1355	Auto	2,163,090	1,012,214
8	1,3-dimethyl-2-nitrobenzene	7:54	474	Auto	1,249,077	607,405
9	triphenyl phosphate (S)	22:01	1321	Auto	1,316,075	652,365
10	perylene d-12 (SURR)	26:47	1607	Auto	904,221	178,904
11	hexachlorocyclopentadiene	10:04	604	Auto	4,314	2,455
12	bis(2-ethylhexyl)adipate	22:01	1321	Auto	61,241	10,447
13	bis(2-ethylhexyl)phthalate	22:53	1373	Auto	1,345,390	442,962
14	2,6-dinitrotoluene	11:52	712	Auto	166,879	80,615
15	2,4-dinitrotoluene	13:00	780	Auto	44,121	7,056
16	butachlor	20:19	1219	Auto	3,443	1,953
17	4,4'-dichlorodiphenyl ether	20:42	1242	Auto	894	894

reversed IS/SURR's

Quantitation Report

Quanfile: 26390I

Quan Entries: 17

Comment: BOHLK; METHOD 525; INST 204; 11/23/01; SAMPLE 26390

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	BV	5.000	UG/L
5	acenaphthene d-10(SURR)	A	732	12:12	BB	3.813	UG/L
6	phenanthrene d-10(SURR)	A	1007	16:47	BV	4.111	UG/L
7	chrysene d-12 (SURR)	A	1355	22:35	BV	3.378	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	474	7:54	BV	4.708	UG/L
9	triphenyl phosphate (S)	A	1321	22:01	BV	5.488	UG/L
10	perylene d-12 (SURR)	A	1607	26:47	BB	3.621	UG/L
11	hexachlorocyclopentadiene	A	604	10:04	BB	0.045	UG/L
16	bis(2-ethylhexyl)adipate	A	1321	22:01	MM	0.126	UG/L
17	bis(2-ethylhexyl)phthalate	A	1373	22:53	BB	1.778	UG/L
20	2,6-dinitrotoluene	A	712	11:52	BB	1.286	UG/L
21	2,4-dinitrotoluene	A	780	13:00	MM	0.299	UG/L
27	butachlor	A	1219	20:19	BB	0.015	UG/L
28	4,4'-dichlorodiphenyl ether	A	1242	20:42	BB	0.004	UG/L

Comments for Sample AD26390 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 12-21-2001 Time: 16:43:32

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.