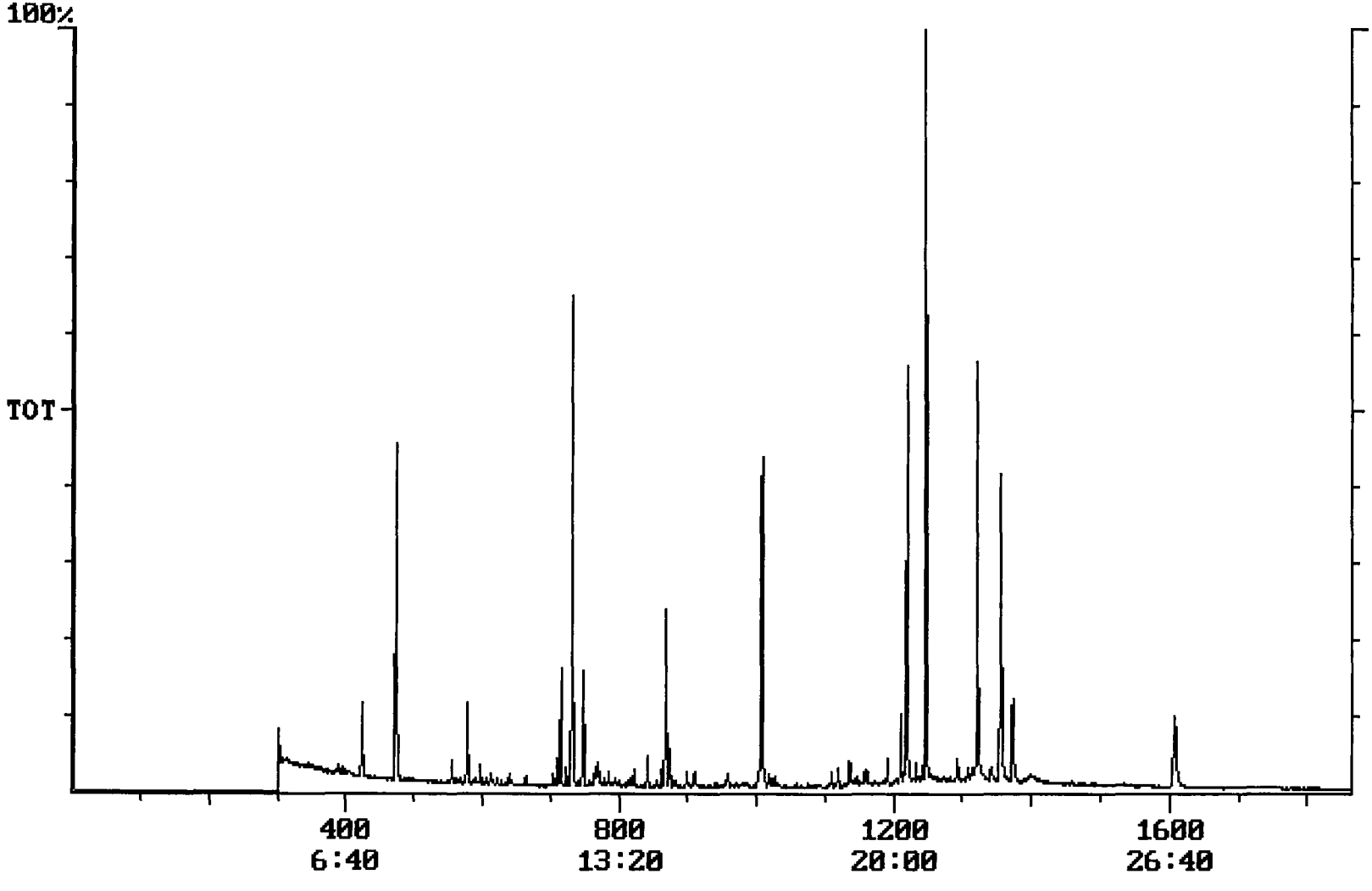
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/20/2001 Time: 15:51:55

Sample: AD26473
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 11/7/01 15:50 Ended: 11/24/01 15:50 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		1.1	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		5.4	1.0
20	Surr_Triphenyl phosphate		5.6	1.0
21	Surr_Perylene-d12		4.2	1.0

Chromatogram Plot File: C:\26473D Date: Nov-24-1991 22:54:19
Comment: BOHLK; METHOD 525; INST 204; 11/24/01; SAMPLE 26473
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 7334560



Integration Report Quanfile: 26473D Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/24/01; SAMPLE 26473
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	1,780,397	1,202,573
2	phenanthrene d-10 (IS)	16:48	1008	Auto	2,870,127	1,009,822
3	chrysene d-12(IS)	22:35	1355	Auto	1,857,362	871,871
4	p-terphenyl d-14(IS)	20:46	1246	Auto	2,972,801	2,205,693
5	acenaphthene d-10(SURR	12:12	732	Auto	1,780,397	1,202,573
6	phenanthrene d-10(SURR	16:48	1008	Auto	2,870,127	1,009,822
7	chrysene d-12 (SURR)	22:35	1355	Auto	1,857,362	871,871
8	1,3-dimethyl-2-nitrobe	7:54	474	Auto	1,145,201	533,430
9	triphenyl phosphate (S	22:01	1321	Auto	1,152,679	509,891
10	perylene d-12 (SURR)	26:47	1607	Auto	905,229	176,392
11	bis(2-ethylhexyl)adipa	22:01	1321	Auto	30,115	6,055
12	bis(2-ethylhexyl)phtha	22:53	1373	Auto	741,499	249,530
13	2,4-dinitrotoluene	13:00	780	Auto	13,067	3,838
14	butachlor	20:19	1219	Auto	2,364	2,364

reviewed FS/SURR's

Quantitation Report Quanfile: 26473D Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/24/01; SAMPLE 26473
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	BV	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1008	16:48	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	BV	3.950	UG/L
6	phenanthrene d-10(SURR	A	1008	16:48	BV	4.432	UG/L
7	chrysene d-12 (SURR)	A	1355	22:35	BV	3.732	UG/L
8	1,3-dimethyl-2-nitrobe	A	474	7:54	BB	5.361	UG/L
9	triphenyl phosphate (S	A	1321	22:01	BV	5.598	UG/L
10	perylene d-12 (SURR)	A	1607	26:47	BB	4.222	UG/L
16	bis(2-ethylhexyl)adipa	A	1321	22:01	MM	0.072	UG/L
17	bis(2-ethylhexyl)phtha	A	1373	22:53	BB	1.111	UG/L
21	2,4-dinitrotoluene	A	780	13:00	BB	0.111	UG/L
27	butachlor	A	1219	20:19	BB	0.012	UG/L

Comments for Sample AD26473 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-22-2001 Time: 11:51:39

The breakdown for Endrin in the Breakdown Standard was 39%. Recoveries for 4 of the 18 compounds in the MDL Standard were outside of their acceptance ranges. Recoveries for 2 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.