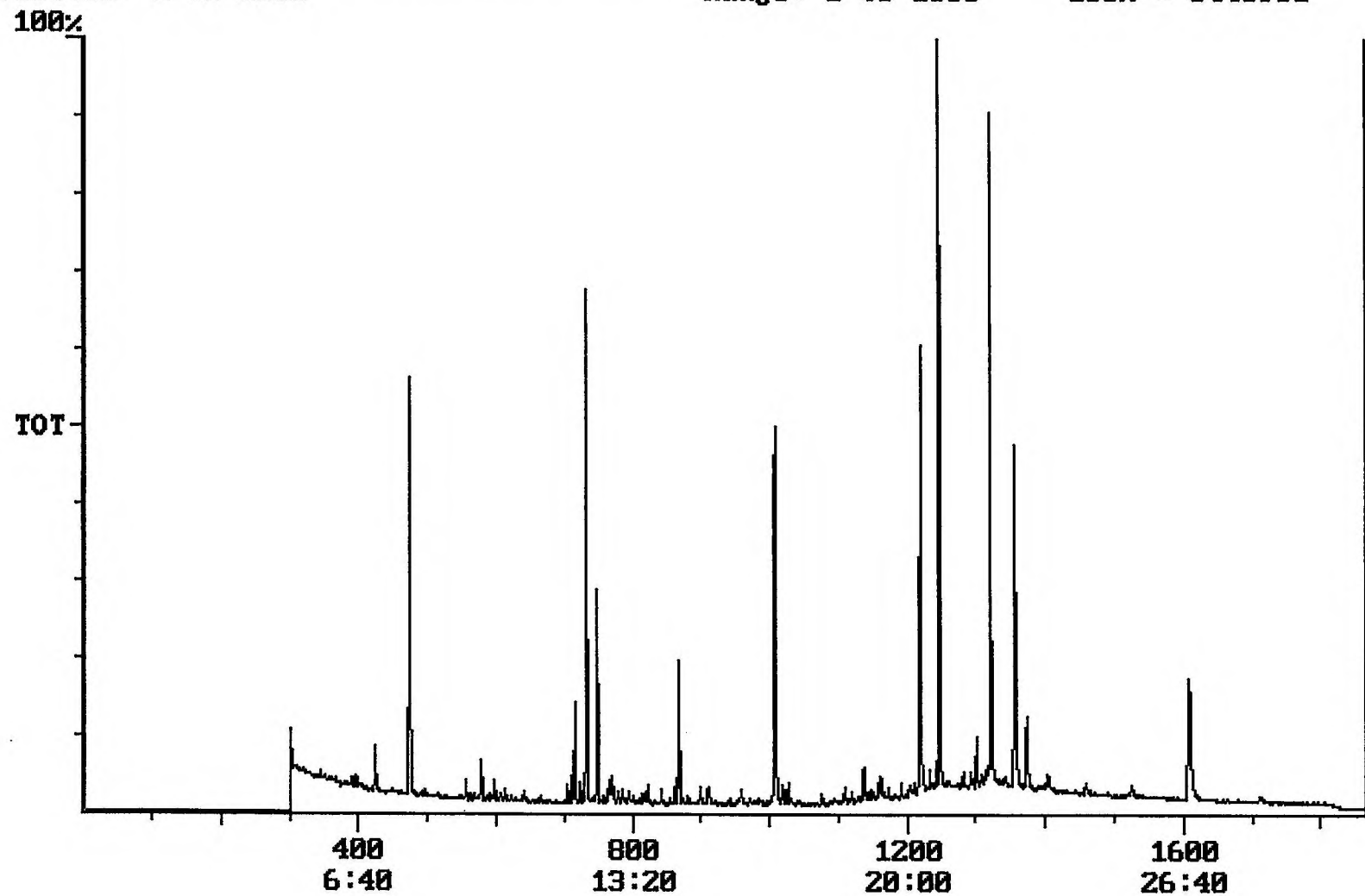


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/21/2001 Time: 16:17:53

Sample: AD26680
Analysis: \$525 Organic Chemicals - 525.2
Units: ug
Started: 11/8/01 16:16 Ended: 11/28/01 16:16 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		0.62	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		5.1	1.0
20	Surr_Triphenyl phosphate		5.8	1.0
21	Surr_Perylene-d12		4.8	1.0

Chromatogram Plot File: E:\26680K Date: Nov-28-1991 22:04:30
Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26680
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 5445931



Integration Report Quanfile: 26680K Quan Entries: 19
 Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26680
 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	1,449,659	844,099
2	phenanthrene d-10 (IS)	16:48	1008	Auto	2,398,470	894,277
3	chrysene d-12(IS)	22:35	1355	Auto	1,778,555	753,748
4	p-terphenyl d-14(IS)	20:46	1246	Auto	2,557,287	1,664,734
5	acenaphthene d-10(SURR)	12:12	732	Auto	1,449,659	844,099
6	phenanthrene d-10(SURR)	16:48	1008	Auto	2,398,470	894,277
7	chrysene d-12 (SURR)	22:35	1355	Auto	1,778,555	753,748
8	1,3-dimethyl-2-nitrobenzene	7:54	474	Auto	891,397	485,203
9	triphenyl phosphate (S)	22:01	1321	Auto	1,150,029	619,402
10	perylene d-12 (SURR)	26:48	1608	Auto	982,472	193,056
11	hexachlorocyclopentadiene	10:04	604	Auto	2,178	1,262
12	hexachlorobenzene	15:14	914	Auto	11,419	4,259
13	simazine	16:10	970	Auto	0	0
14	bis(2-ethylhexyl)adipate	21:54	1314	Auto	27,258	6,553
15	bis(2-ethylhexyl)phthalate	22:53	1373	Auto	407,355	143,108
16	2,6-dinitrotoluene	11:52	712	Auto	98,142	43,458
17	2,4-dinitrotoluene	13:00	780	Auto	9,236	3,568
18	butachlor	20:19	1219	Auto	2,511	2,511
19	4,4'-dichlorodiphenyl ether	20:42	1242	Auto	8,129	5,481

reviewed IS/SURR's

Quantitation Report Quanfile: 26680K Quan Entries: 19
 Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26680
 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	1008	16:48	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1355	22:35	BB	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	3.739	UG/L
6	phenanthrene d-10(SURR	A	1008	16:48	BV	4.306	UG/L
7	chrysene d-12 (SURR)	A	1355	22:35	BB	4.154	UG/L
8	1,3-dimethyl-2-nitroben	A	474	7:54	BV	5.125	UG/L
9	triphenyl phosphate (S	A	1321	22:01	BB	5.833	UG/L
10	perylene d-12 (SURR)	A	1608	26:48	BB	4.785	UG/L
11	hexachlorocyclopentadi	A	604	10:04	BB	0.035	UG/L
12	hexachlorobenzene	A	914	15:14	BB	0.066	UG/L
13	simazine	A	970	16:10	BB	0.001	UG/L
16	bis(2-ethylhexyl)adipa	A	1314	21:54	MM	0.068	UG/L
17	bis(2-ethylhexyl)phtha	A	1373	22:53	BV	0.624	UG/L
20	2,6-dinitrotoluene	A	712	11:52	BB	1.154	UG/L
21	2,4-dinitrotoluene	A	780	13:00	BB	0.096	UG/L
27	butachlor	A	1219	20:19	BB	0.015	UG/L
28	4,4'-dde	A	1242	20:42	BB	0.042	UG/L

Comments for Sample AD26680 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 12-22-2001 Time: 12:59:38

The breakdown for Endrin in the Breakdown Standard was 40%. Recoveries for 4 of the 18 compounds in the MDL Standard were outside of their acceptance ranges. Recoveries for 4 of the 18 compounds in the LCS Standard were below their acceptance ranges.