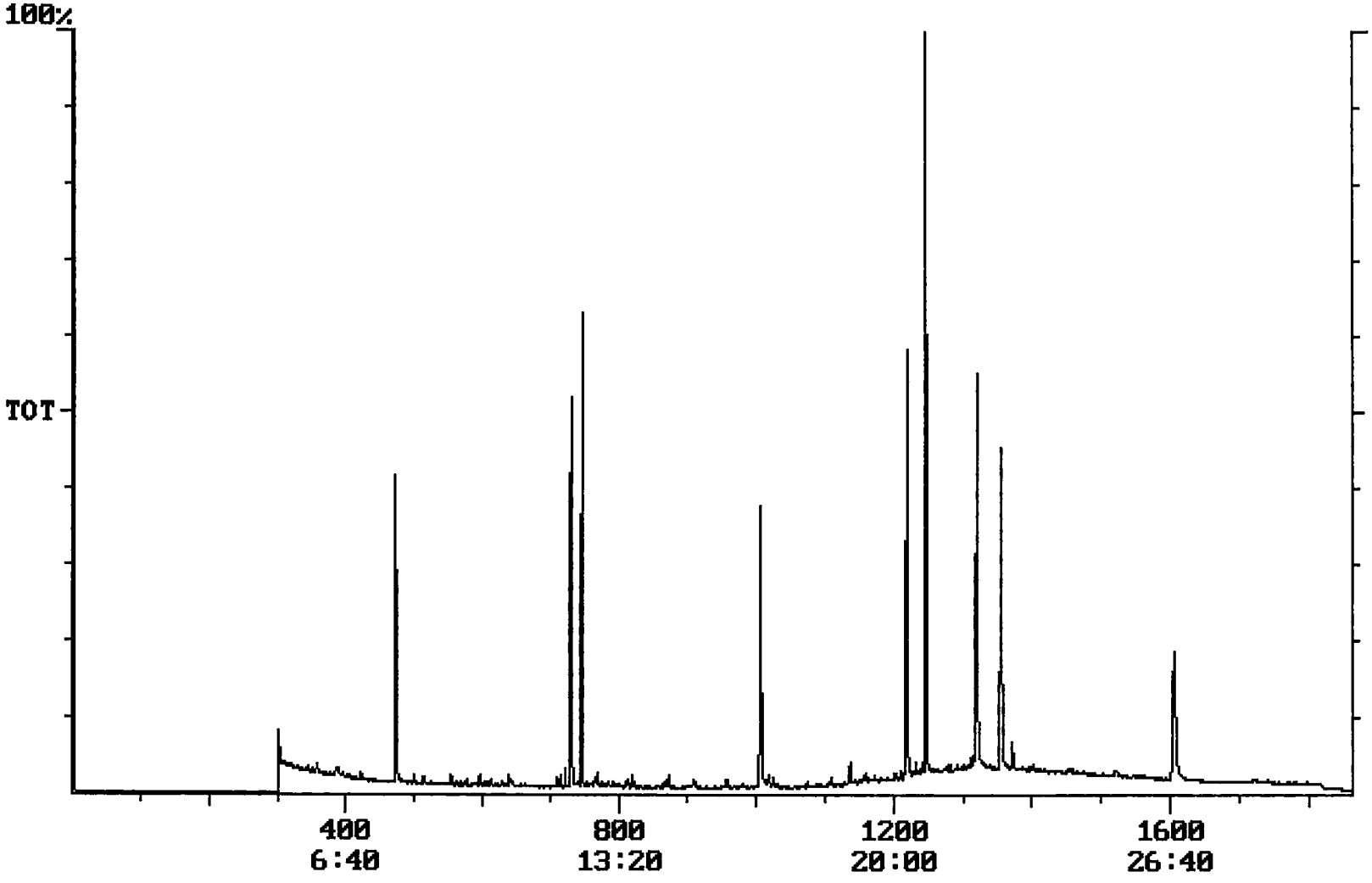
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 01/08/2002 Time: 13:18:02

Sample: AD27625
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 11/23/01 13:14 Ended: 12/8/01 13:14 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		< MDL		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		5.3		1.0
20	Surr_Triphenyl phosphate		6.0		1.0
21	Surr_Perylene-d12		4.9		1.0

Chromatogram Plot File: E:\27625DD Date: Dec-08-1991 22:00:33
Comment: BOHLK; METHOD 525; INST 204; 12/08/01; SAMPLE 27625
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1859 Range: 1 to 1859 100% = 9866002



Integration Report Quanfile: 27625DD Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 12/08/01; SAMPLE 27625
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:11	731	Auto	2,426,671	1,324,567
2	phenanthrene d-10 (IS)	16:46	1006	Auto	3,557,238	1,219,263
3	chrysene d-12(IS)	22:34	1354	Auto	2,792,025	1,296,583
4	p-terphenyl d-14(IS)	20:45	1245	Auto	4,209,381	3,199,133
5	acenaphthene d-10(SURR	12:11	731	Auto	2,426,671	1,324,567
6	phenanthrene d-10(SURR	16:46	1006	Auto	3,557,238	1,219,263
7	chrysene d-12 (SURR)	22:34	1354	Auto	2,792,025	1,296,583
8	1,3-dimethyl-2-nitrobe	7:52	472	Auto	1,199,873	658,272
9	triphenyl phosphate (S	21:59	1319	Auto	1,622,109	768,421
10	perylene d-12 (SURR)	26:45	1605	Auto	1,989,727	440,109
11	bis(2-ethylhexyl)adipa	21:52	1312	Auto	15,640	5,524
12	bis(2-ethylhexyl)phtha	22:51	1371	Auto	231,475	104,511
13	benzo(a)pyrene	26:33	1593	Auto	0	0
14	2,6-dinitrotoluene	11:51	711	Auto	25,597	16,251
15	butachlor	20:18	1218	Auto	15,677	7,964

reviewed IS/surr's

Quantitation Report Quanfile: 27625DD Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 12/08/01; SAMPLE 27625
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	731	12:11	BV	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1006	16:46	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1354	22:34	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1245	20:45	BV	5.000	UG/L
5	acenaphthene d-10(SURR	A	731	12:11	BV	3.175	UG/L
6	phenanthrene d-10(SURR	A	1006	16:46	BV	3.635	UG/L
7	chrysene d-12 (SURR)	A	1354	22:34	BB	3.487	UG/L
8	1,3-dimethyl-2-nitrobe	A	472	7:52	BV	5.299	UG/L
9	triphenyl phosphate (S	A	1319	21:59	BB	5.996	UG/L
10	perylene d-12 (SURR)	A	1605	26:45	BB	4.891	UG/L
16	bis(2-ethylhexyl)adipa	A	1312	21:52	MM	0.038	UG/L
17	bis(2-ethylhexyl)phtha	A	1371	22:51	VV	0.283	UG/L
18	benzo(a)pyrene	A	1593	26:33	MM	0.001	UG/L
20	2,6-dinitrotoluene	A	711	11:51	BB	0.179	UG/L
27	butachlor	A	1218	20:18	BB	0.049	UG/L

Comments for Sample AD27625 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 01-22-2002 Time: 14:17:25

The recoveries for 2 of the 18 compounds in the LCS Standard were below their acceptance ranges. The recoveries for 3 of the 18 compounds in the Spiked Sample were outside of their acceptance ranges. The breakdown for Endrin in the Breakdown Standard was 61%. The injection port is kept at 280 degrees Celcius which may be contributing to high Endrin Breakdown. An investigation into modifying injection port parameters will be made.