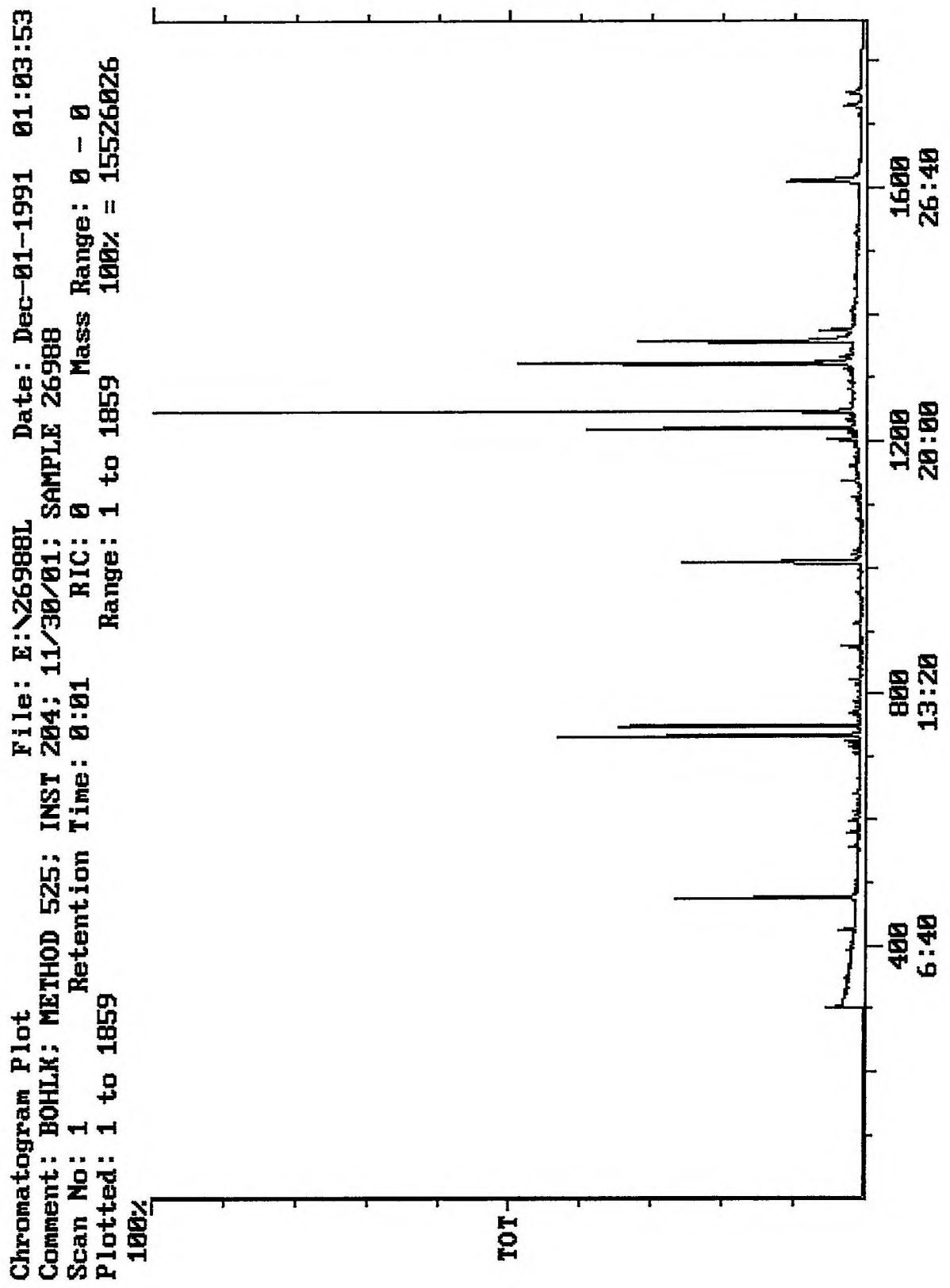
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 01/03/2002 Time: 12:18:19

Sample: AD26988
Analysis: S525 Organic Chemicals - 525.2
Units: ug
Started: 11/13/01 12:17 Ended: 11/30/01 12:17 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-nitrobenze		4.2		1.0
20	Surr_Triphenyl phosphate		5.1		1.0
21	Surr_Perylene-d12		5.4		1.0



Integration Report Quanfile: 26988L Quan Entries: 15
Comment: BOHLK: METHOD 525; INST 204; 11/30/01; SAMPLE 26988
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:13	733	Auto	2,633,987	1,830,607
2	phenanthrene d-10 (IS)	16:49	1009	Auto	3,572,801	1,377,091
3	chrysene d-12(IS)	22:36	1356	Auto	3,106,795	1,305,014
4	p-terphenyl d-14(IS)	20:47	1247	Auto	7,176,621	5,470,000
5	acenaphthene d-10(SURR	12:13	733	Auto	2,633,987	1,830,607
6	phenanthrene d-10(SURR	16:49	1009	Auto	3,572,801	1,377,091
7	chrysene d-12 (SURR)	22:36	1356	Auto	3,106,795	1,305,014
8	1,3-dimethyl-2-nitrobe	7:55	475	Auto	1,316,101	670,907
9	triphenyl phosphate (S	22:02	1322	Auto	1,753,780	989,754
10	perylene d-12 (SURR)	26:50	1610	Auto	1,953,148	433,619
11	hexachlorocyclopentadi	10:04	604	Auto	2,777	1,433
12	bis(2-ethylhexyl)adipa	22:02	1322	Auto	46,446	7,739
13	bis(2-ethylhexyl)phtha	22:54	1374	Auto	565,697	203,170
14	2,4-dinitrotoluene	13:02	782	Auto	1,148	1,148
15	butachlor	20:20	1220	Auto	3,267	2,102

reviewed IS/ surr's

Quantitation Report Quanfile: 26988L Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 11/30/01; SAMPLE 26988
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	733	12:13	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1009	16:49	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1356	22:36	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1247	20:47	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	733	12:13	BB	2.421	UG/L
6	phenanthrene d-10(SURR	A	1009	16:49	BV	2.286	UG/L
7	chrysene d-12 (SURR)	A	1356	22:36	BB	2.586	UG/L
8	1,3-dimethyl-2-nitrobe	A	475	7:55	BV	4.165	UG/L
9	triphenyl phosphate (S	A	1322	22:02	BV	5.092	UG/L
10	perylene d-12 (SURR)	A	1610	26:50	BB	5.446	UG/L
11	hexachlorocyclopentadi	A	604	10:04	BB	0.024	UG/L
16	bis(2-ethylhexyl)adipa	A	1322	22:02	MM	0.066	UG/L
17	bis(2-ethylhexyl)phtha	A	1374	22:54	BV	0.493	UG/L
21	2,4-dinitrotoluene	A	782	13:02	BB	0.007	UG/L
27	butachlor	A	1220	20:20	BB	0.013	UG/L

Comments for Sample AD26988 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 01-04-2002 Time: 10:43:39

The recoveries for 5 of the 18 compounds in the LCS Standard were outside of 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 48%.