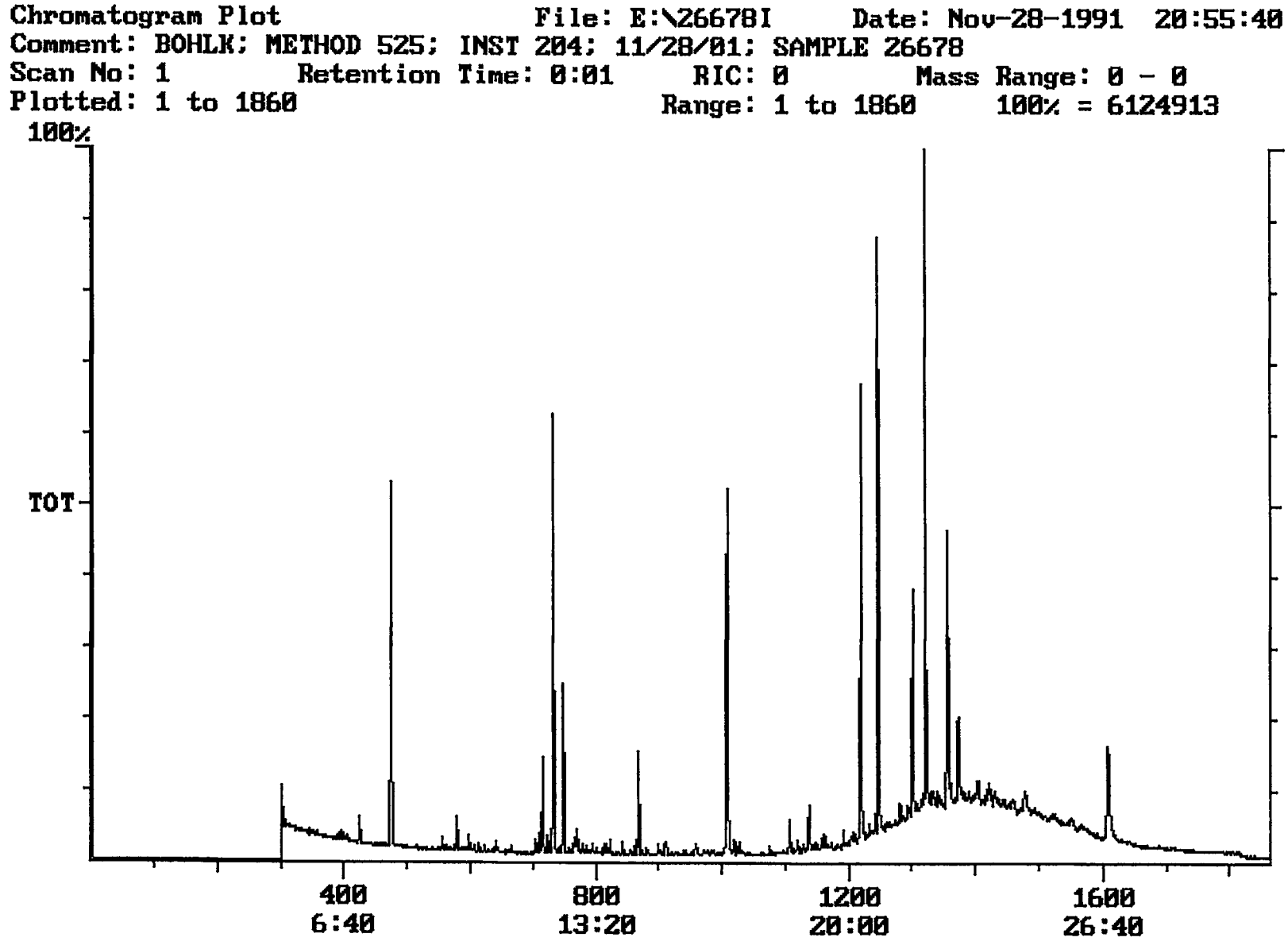
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/21/2001 Time: 16:15:00

Sample: AD26678
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 11/8/01 16:13 Ended: 11/28/01 16:13 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		0.69	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		5.2	1.0
20	Surr_Triphenyl phosphate	USPEC	6.9	1.0
21	Surr_Perylene-d12		4.5	1.0



Integration Report Quanfile: 26678I Quan Entries: 17
Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26678
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	1,637,360	902,662
2	phenanthrene d-10 (IS)	16:48	1008	Auto	2,587,482	1,023,217
3	chrysene d-12(IS)	22:35	1355	Auto	1,798,898	722,118
4	p-terphenyl d-14(IS)	20:46	1246	Auto	2,677,768	1,613,372
5	acenaphthene d-10(SURR	12:12	732	Auto	1,637,360	902,662
6	phenanthrene d-10(SURR	16:48	1008	Auto	2,587,482	1,023,217
7	chrysene d-12 (SURR)	22:35	1355	Auto	1,798,898	722,118
8	1,3-dimethyl-2-nitrobe	7:54	474	Auto	1,016,838	605,870
9	triphenyl phosphate (S	22:01	1321	Auto	1,368,519	763,971
10	perylene d-12 (SURR)	26:48	1608	Auto	940,499	195,726
11	hexachlorocyclopentadi	10:04	604	Auto	1,512	1,512
12	hexachlorobenzene	15:13	913	Auto	11,355	4,725
13	simazine	16:04	964	Auto	0	0
14	bis(2-ethylhexyl)phtha	22:53	1373	Auto	456,835	172,304
15	2,4-dinitrotoluene	13:01	781	Auto	7,997	2,811
16	butachlor	20:22	1222	Auto	0	0
17	4,4'-dde	20:42	1242	Auto	11,262	7,652

reviewed ES/ Surr's

Quantitation Report Quanfile: 26678I Quan Entries: 17
Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26678
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	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	BB	4.033	UG/L
6	phenanthrene d-10(SURR	A	1008	16:48	BV	4.436	UG/L
7	chrysene d-12 (SURR)	A	1355	22:35	BV	4.013	UG/L
8	1,3-dimethyl-2-nitrobe	A	474	7:54	BV	5.176	UG/L
9	triphenyl phosphate (S	A	1321	22:01	BV	6.862	UG/L
10	perylene d-12 (SURR)	A	1608	26:48	BB	4.529	UG/L
11	hexachlorocyclopentadi	A	604	10:04	BB	0.021	UG/L
12	hexachlorobenzene	A	913	15:13	BB	0.058	UG/L
13	simazine	A	964	16:04	BB	0.001	UG/L
17	bis(2-ethylhexyl)phtha	A	1373	22:53	VB	0.694	UG/L
21	2,4-dinitrotoluene	A	781	13:01	BB	0.074	UG/L
27	butachlor	A	1222	20:22	BB	0.001	UG/L
28	4,4'-dde	A	1242	20:42	BB	0.053	UG/L

Comments for Sample AD26678 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-22-2001 Time: 12:57:03

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. The recovery for 1 of the 3 Surrogate Standards in this sample was above its acceptance range.