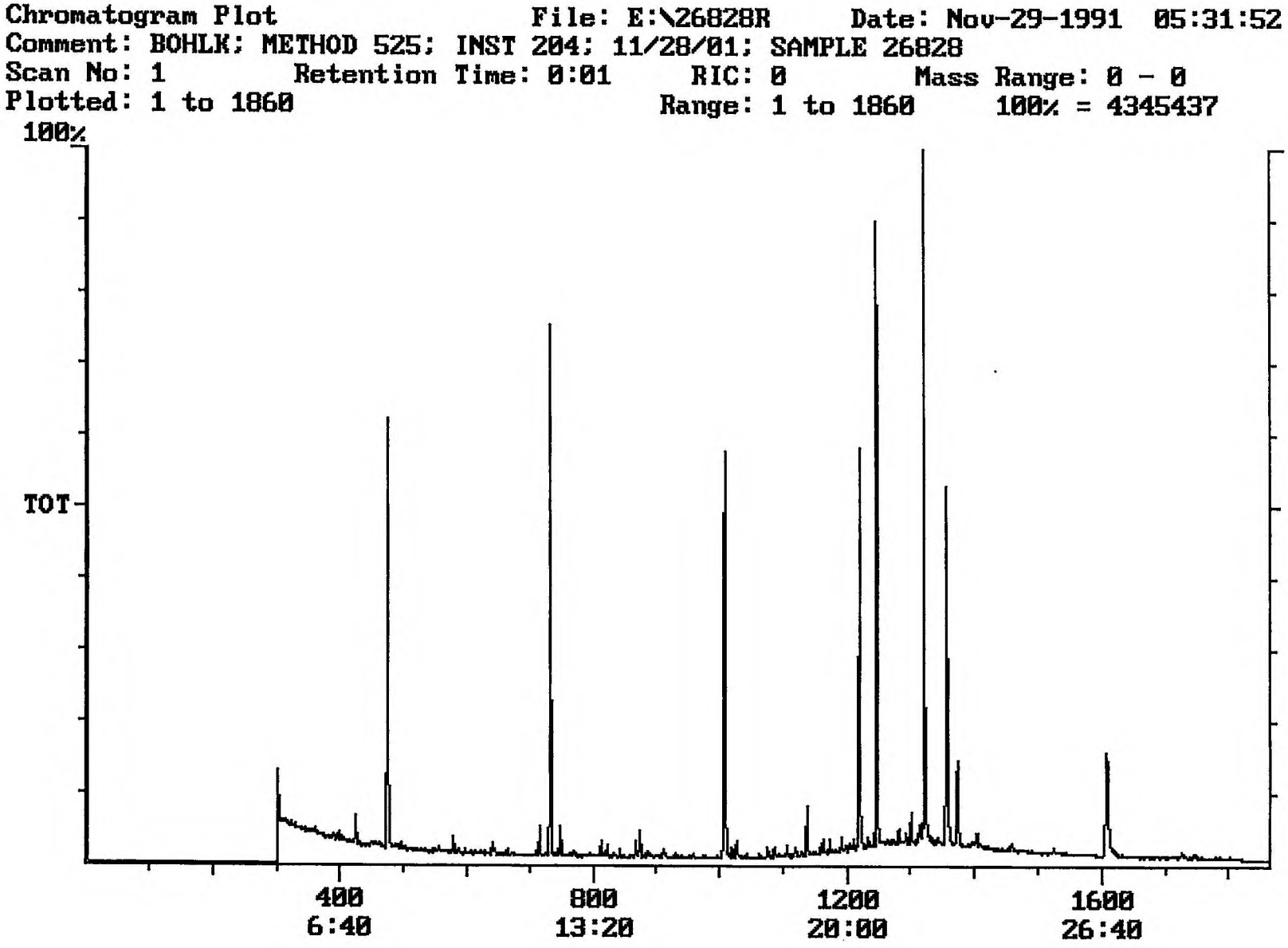
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/21/2001 Time: 16:27:23

Sample: AD26828
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 11/8/01 16:26 Ended: 11/28/01 16:26 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		0.74	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		4.8	1.0
20	Surr_Triphenyl phosphate		6.1	1.0
21	Surr_Perylene-d12		4.6	1.0



Integration Report Quanfile: 26828R Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26828
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	1,288,009	763,349
2	phenanthrene d-10 (IS)	16:48	1008	Auto	2,048,285	806,058
3	chrysene d-12(IS)	22:35	1355	Auto	1,531,877	665,209
4	p-terphenyl d-14(IS)	20:46	1246	Auto	2,196,941	1,191,451
5	acenaphthene d-10(SURR	12:12	732	Auto	1,288,009	763,349
6	phenanthrene d-10(SURR	16:48	1008	Auto	2,048,285	806,058
7	chrysene d-12 (SURR)	22:35	1355	Auto	1,531,877	665,209
8	1,3-dimethyl-2-nitrobe	7:54	474	Auto	743,488	413,317
9	triphenyl phosphate (S	22:01	1321	Auto	1,036,548	610,500
10	perylene d-12 (SURR)	26:48	1608	Auto	811,398	164,164
11	bis(2-ethylhexyl)adipa	21:54	1314	Auto	28,349	6,213
12	bis(2-ethylhexyl)phtha	22:53	1373	Auto	413,078	150,829
13	2,6-dinitrotoluene	11:52	712	Auto	16,901	7,817
14	butachlor	20:20	1220	Auto	1,071	1,071

reviewed IS/surr's

Quantitation Report Quanfile: 26828R Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26828
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	3.867	UG/L
6	phenanthrene d-10(SURR	A	1008	16:48	BV	4.280	UG/L
7	chrysene d-12 (SURR)	A	1355	22:35	BV	4.165	UG/L
8	1,3-dimethyl-2-nitrobe	A	474	7:54	BB	4.811	UG/L
9	triphenyl phosphate (S	A	1321	22:01	BV	6.104	UG/L
10	perylene d-12 (SURR)	A	1608	26:48	BB	4.588	UG/L
16	bis(2-ethylhexyl)adipa	A	1314	21:54	MM	0.082	UG/L
17	bis(2-ethylhexyl)phtha	A	1373	22:53	BV	0.738	UG/L
20	2,6-dinitrotoluene	A	712	11:52	BB	0.225	UG/L
27	butachlor	A	1220	20:20	BB	0.008	UG/L

Comments for Sample AD26828 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-22-2001 Time: 13:07:05

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.