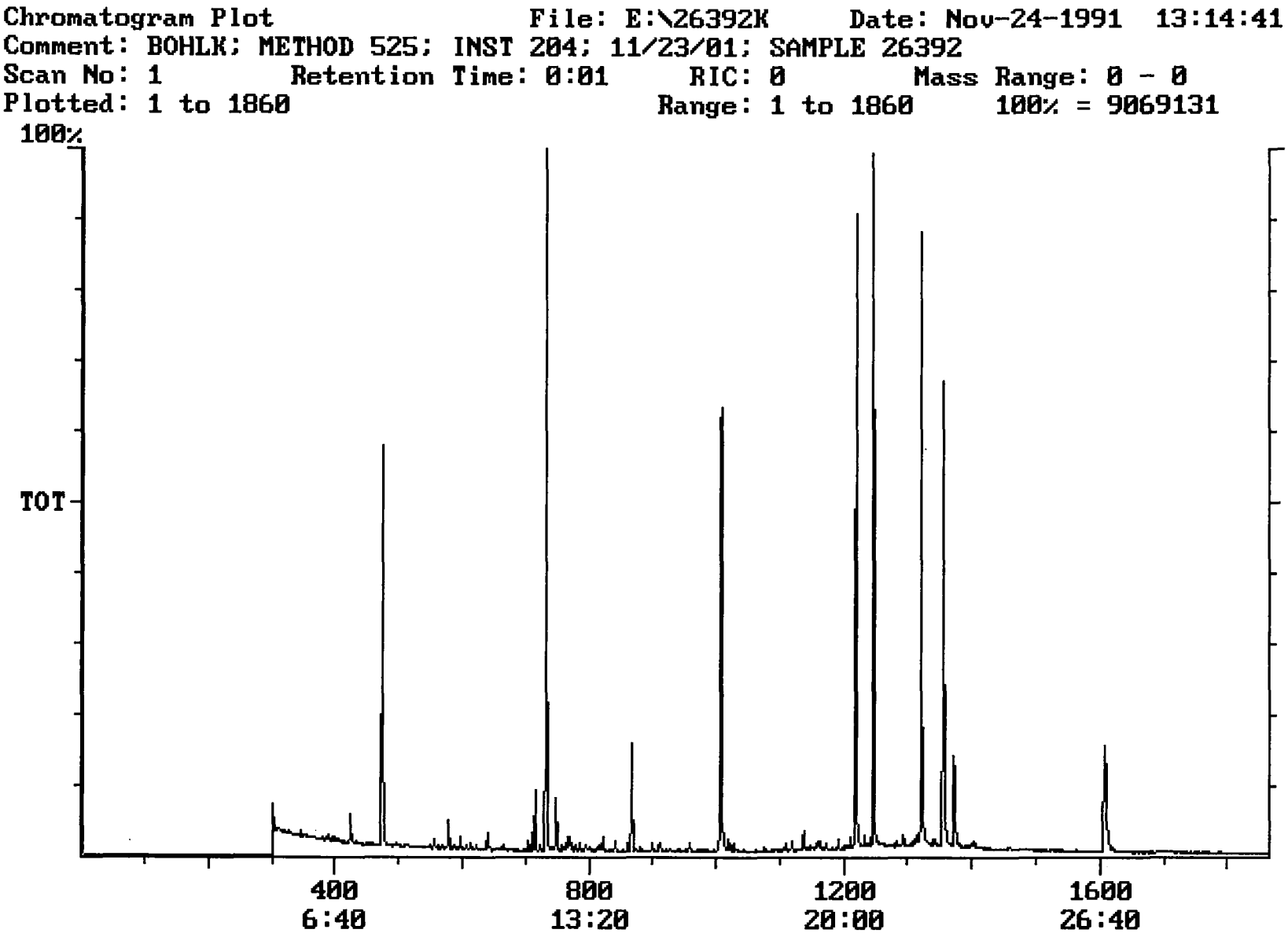
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/20/2001 Time: 09:34:50

Sample: AD26392
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 11/6/01 09:33 Ended: 11/23/01 09:33 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		0.68	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		4.2	1.0
20	Surr_Triphenyl phosphate		5.0	1.0
21	Surr_Perylene-d12		3.7	1.0



Integration Report Quanfile: 26392K Quan Entries: 17
Comment: BOHLK; METHOD 525; INST 204; 11/23/01; SAMPLE 26392
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	3,696,810	2,445,890
2	phenanthrene d-10 (IS)	16:47	1007	Auto	5,565,045	1,989,967
3	chrysene d-12(IS)	22:35	1355	Auto	4,042,778	1,923,705
4	p-terphenyl d-14(IS)	20:46	1246	Auto	3,909,170	2,890,789
5	acenaphthene d-10(SURR	12:12	732	Auto	3,696,810	2,445,890
6	phenanthrene d-10(SURR	16:47	1007	Auto	5,565,045	1,989,967
7	chrysene d-12 (SURR)	22:35	1355	Auto	4,042,778	1,923,705
8	1,3-dimethyl-2-nitrobe	7:54	474	Auto	1,851,091	989,044
9	triphenyl phosphate (S	22:01	1321	Auto	2,223,680	1,059,442
10	perylene d-12 (SURR)	26:47	1607	Auto	1,712,876	354,921
11	hexachlorocyclopentadi	10:03	603	Auto	1,082	1,082
12	bis(2-ethylhexyl)adipa	22:01	1321	Auto	42,988	8,950
13	bis(2-ethylhexyl)phtha	22:52	1372	Auto	997,137	349,994
14	benzo(a)pyrene	26:48	1608	Auto	6,126	2,962
15	2,6-dinitrotoluene	11:52	712	Auto	95,898	50,000
16	2,4-dinitrotoluene	13:00	780	Auto	15,096	3,517
17	butachlor	20:18	1218	Auto	865	865

reviewed IS/ surr's

Quantitation Report Quanfile: 26392K Quan Entries: 17
Comment: BOHLK; METHOD 525; INST 204; 11/23/01; SAMPLE 26392
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	1007	16:47	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	BV	5.000	UG/L
5	acenaphthene d-10(SURR	A	732	12:12	BB	6.237	UG/L
6	phenanthrene d-10(SURR	A	1007	16:47	BV	6.535	UG/L
7	chrysene d-12 (SURR)	A	1355	22:35	BV	6.177	UG/L
8	1,3-dimethyl-2-nitrobe	A	474	7:54	BB	4.173	UG/L
9	triphenyl phosphate (S	A	1321	22:01	BB	4.962	UG/L
10	perylene d-12 (SURR)	A	1607	26:47	BB	3.670	UG/L
11	hexachlorocyclopentadi	A	603	10:03	BB	0.007	UG/L
16	bis(2-ethylhexyl)adipa	A	1321	22:01	MM	0.047	UG/L
17	bis(2-ethylhexyl)phtha	A	1372	22:52	BB	0.674	UG/L
18	benzo(a)pyrene	A	1608	26:48	BB	0.007	UG/L
20	2,6-dinitrotoluene	A	712	11:52	BB	0.444	UG/L
21	2,4-dinitrotoluene	A	780	13:00	BB	0.062	UG/L
27	butachlor	A	1218	20:18	BB	0.003	UG/L

Comments for Sample AD26392 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-21-2001 Time: 16:46:04

The breakdown for Endrin in the Breakdown Standard was 56%. Recoveries for 4 of the 18 compounds in the MDL Standard were outside of their acceptance ranges. Recoveries for 3 of the 18 compounds in the LCS Standard were outside of their acceptance ranges. Recoveries for 3 of the 18 compounds in the Spiked Sample were outside of their acceptance ranges. The level for bis(2-Ethylhexyl)phthalate was 0.88 ug/L in the Laboratory Blank.