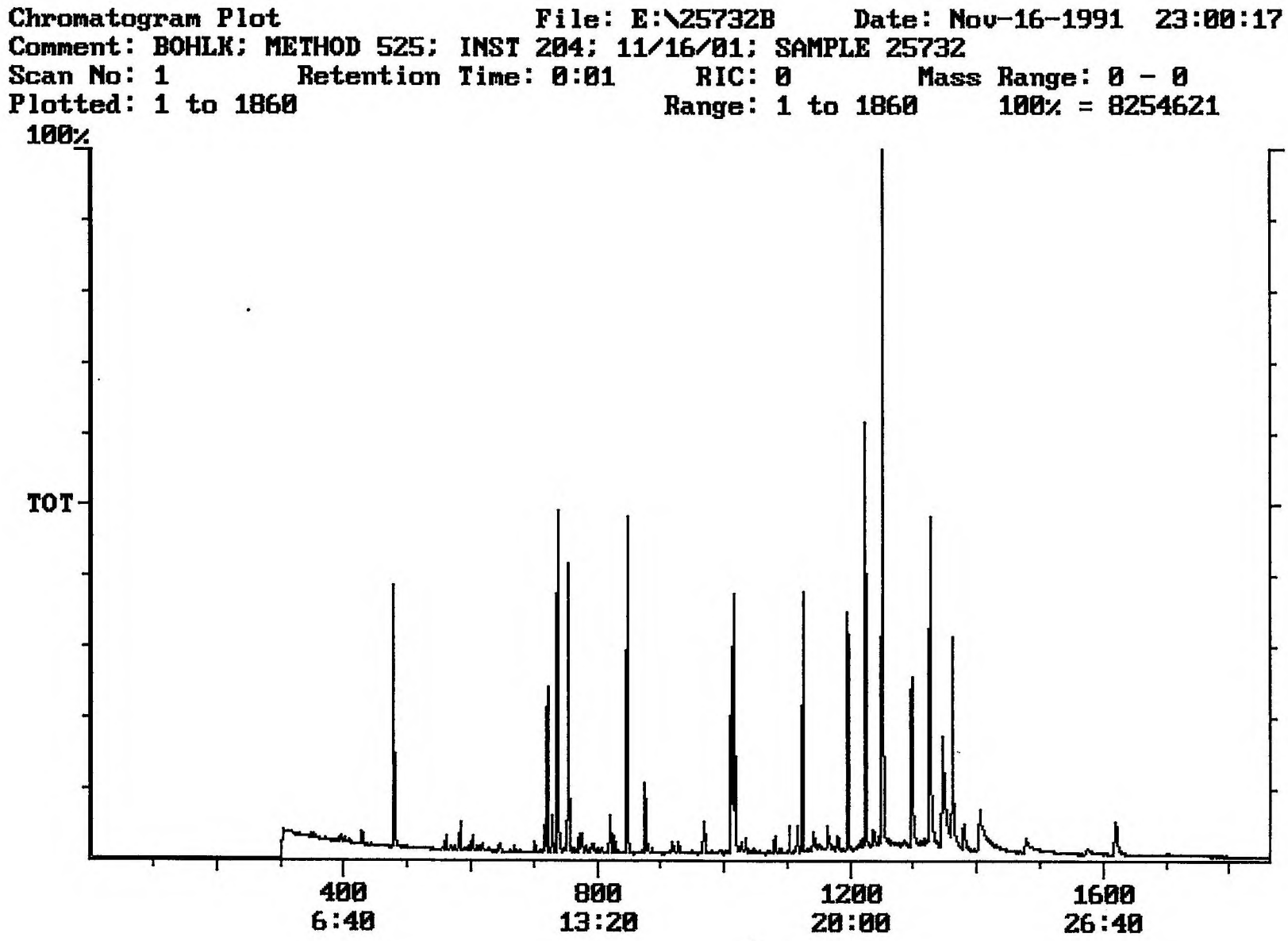
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 11/30/2001 Time: 09:32:35

Sample: AD25732
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 11/1/01 09:28 Ended: 11/16/01 09:28 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		5.2		1.0
20	Surr_Triphenyl phosphate		5.5		1.0
21	Surr_Perylene-d12		2.9		1.0



Integration Report Quanfile: 25732B Quan Entries: 12
Comment: BOHLK; METHOD 525; INST 204; 11/16/01; SAMPLE 25732
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:18	738	Auto	1,570,713	908,333
2	phenanthrene d-10 (IS)	16:56	1016	Auto	2,586,189	1,009,362
3	chrysene d-12(IS)	22:41	1361	Auto	1,669,019	700,321
4	p-terphenyl d-14(IS)	20:51	1251	Auto	3,150,155	1,706,001
5	acenaphthene d-10(SURR	12:18	738	Auto	1,570,713	908,333
6	phenanthrene d-10(SURR	16:56	1016	Auto	2,586,189	1,009,362
7	chrysene d-12 (SURR)	22:41	1361	Auto	1,669,019	700,321
8	1,3-dimethyl-2-nitrobe	7:58	478	Auto	1,010,527	484,469
9	triphenyl phosphate (S	22:06	1326	Auto	1,305,566	555,438
10	perylene d-12 (SURR)	27:00	1620	Auto	582,541	90,060
11	bis(2-ethylhexyl)phtha	22:59	1379	Auto	319,781	79,313
12	2,6-dinitrotoluene	11:58	718	Auto	84,286	39,589

Reviewed 15/Sm

Quantitation Report Quanfile: 25732B Quan Entries: 12
Comment: BOHLK; METHOD 525; INST 204; 11/16/01; SAMPLE 25732
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	738	12:18	BV	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1016	16:56	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1361	22:41	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1251	20:51	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	738	12:18	BV	3.281	UG/L
6	phenanthrene d-10(SURR	A	1016	16:56	BV	3.523	UG/L
7	chrysene d-12 (SURR)	A	1361	22:41	BB	3.409	UG/L
8	1,3-dimethyl-2-nitrobe	A	478	7:58	BB	5.165	UG/L
9	triphenyl phosphate (S	A	1326	22:06	BV	5.456	UG/L
10	perylene d-12 (SURR)	A	1620	27:00	BB	2.892	UG/L
17	bis(2-ethylhexyl)phtha	A	1379	22:59	BB	0.490	UG/L
20	2,6-dinitrotoluene	A	718	11:58	BB	0.837	UG/L

Comments for Sample AD25732 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-11-2001 Time: 11:00:42

Recoveries for 2 of the 18 compounds in the MDL Standard were above thier acceptance ranges. Recoveries for 3 of the 18 compounds in the LCS Standard were outside of their acceptance ranges. Recoveries for 4 of the 18 compounds in the Spiked Sample were outside of their acceptance ranges. The recovery for 1 of the 3 Surrogate Standards in this sample was below its acceptance range.