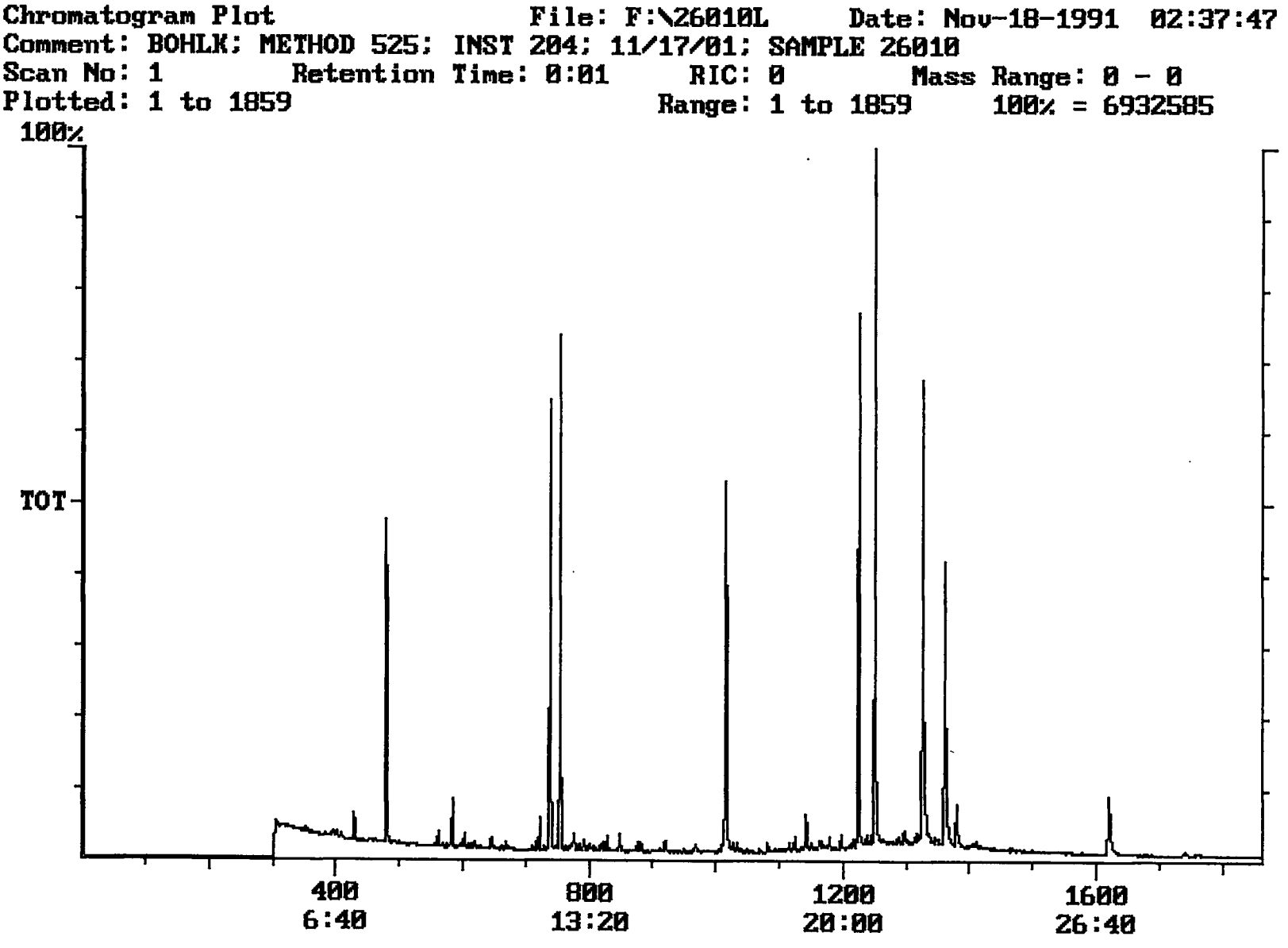
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/13/2001 Time: 18:29:31

Sample: AD26010
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 12/1/01 18:28 Ended: 11/17/01 18:28 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-nitrobenze		5.3		1.0
20	Surr_Triphenyl phosphate		5.4		1.0
21	Surr_Perylene-d12		3.5		1.0



Integration Report Quanfile: 26010L Quan Entries: 16
Comment: BOHLK; METHOD 525; INST 204; 11/17/01; SAMPLE 26010
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	1,818,315	1,013,440
2	phenanthrene d-10 (IS)	16:56	1016	Auto	2,938,994	1,196,013
3	chrysene d-12(IS)	22:41	1361	Auto	1,858,796	823,207
4	p-terphenyl d-14(IS)	20:51	1251	Auto	3,380,593	2,057,924
5	acenaphthene d-10(SURR	12:18	738	Auto	1,818,315	1,013,440
6	phenanthrene d-10(SURR	16:56	1016	Auto	2,938,994	1,196,013
7	chrysene d-12 (SURR)	22:41	1361	Auto	1,858,796	823,207
8	1,3-dimethyl-2-nitrobe	7:58	478	Auto	1,194,608	600,605
9	triphenyl phosphate (S	22:06	1326	Auto	1,448,042	672,467
10	perylene d-12 (SURR)	26:59	1619	Auto	779,439	142,439
11	hexachlorocyclopentadi	10:08	608	Auto	1,238	1,238
12	hexachlorobenzene	15:22	922	Auto	16,435	5,674
13	bis(2-ethylhexyl)adipa	21:59	1319	Auto	41,041	3,857
14	bis(2-ethylhexyl)phtha	22:59	1379	Auto	407,615	110,828
15	2,6-dinitrotoluene	11:58	718	Auto	24,879	12,566
16	4,4'-dde	20:46	1246	Auto	10,241	7,036

renewed IS/ surr's

Quantitation Report Quanfile: 26010L Quan Entries: 16
Comment: BOHLK; METHOD 525; INST 204; 11/17/01; SAMPLE 26010
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	BV	5.000	UG/L
5	acenaphthene d-10(SURR	A	738	12:18	BV	3.539	UG/L
6	phenanthrene d-10(SURR	A	1016	16:56	BV	3.731	UG/L
7	chrysene d-12 (SURR)	A	1361	22:41	BB	3.538	UG/L
8	1,3-dimethyl-2-nitrobe	A	478	7:58	BV	5.275	UG/L
9	triphenyl phosphate (S	A	1326	22:06	BB	5.434	UG/L
10	perylene d-12 (SURR)	A	1619	26:59	BB	3.474	UG/L
11	hexachlorocyclopentadi	A	608	10:08	BB	0.015	UG/L
12	hexachlorobenzene	A	922	15:22	BB	0.066	UG/L
16	bis(2-ethylhexyl)adipa	A	1319	21:59	MM	0.183	UG/L
17	bis(2-ethylhexyl)phtha	A	1379	22:59	BB	0.563	UG/L
20	2,6-dinitrotoluene	A	718	11:58	BB	0.210	UG/L
28	4,4'-dde	A	1246	20:46	BB	0.041	UG/L

Comments for Sample AD26010 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-15-2001 Time: 11:41:03

Recoveries for 3 of the 18 compounds in the MDL and LCS Standards are outside of their acceptance ranges. Recoveries for 4 of the 18 compounds in the Spiked Sample are outside of their acceptance ranges.