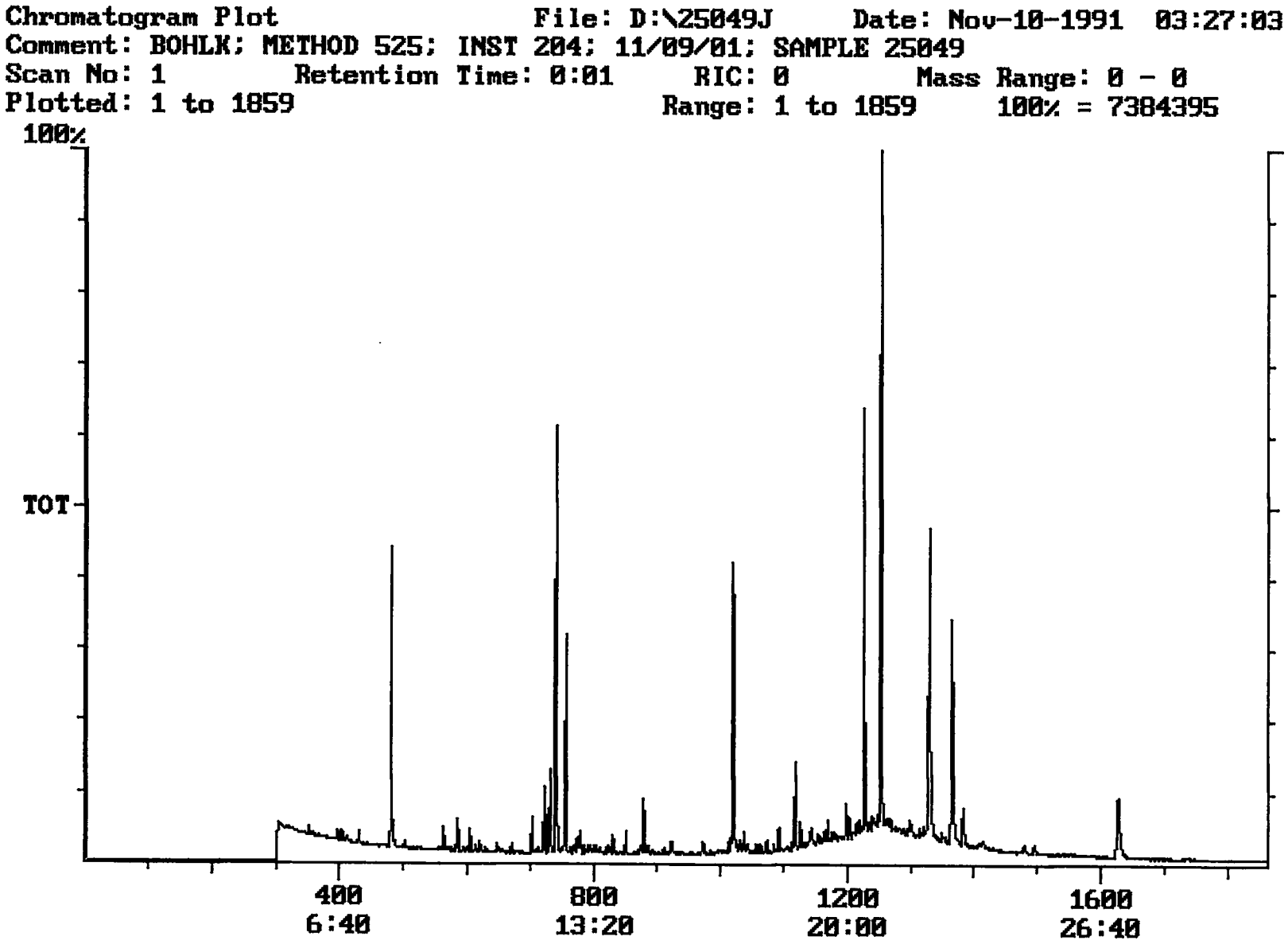
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 11/29/2001 Time: 10:29:17

Sample: AD25049
Analysis: S525 Organic Chemicals - 525.2
Units: ug
Started: 10/25/01 10:04 Ended: 11/9/01 10:04 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		5.0		1.0
20	Surr_Triphenyl phosphate		5.1		1.0
21	Surr_Perylene-d12		3.3		1.0



Integration Report Quanfile: 25049J Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 11/09/01; SAMPLE 25049
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:20	740	Auto	1,643,085	1,012,634
2	phenanthrene d-10 (IS)	16:59	1019	Auto	2,523,091	1,011,089
3	chrysene d-12(IS)	22:44	1364	Auto	1,919,676	690,512
4	p-terphenyl d-14(IS)	20:53	1253	Auto	3,417,569	2,075,247
5	acenaphthene d-10(SURR	12:20	740	Auto	1,643,085	1,012,634
6	phenanthrene d-10(SURR	16:59	1019	Auto	2,523,091	1,011,089
7	chrysene d-12 (SURR)	22:44	1364	Auto	1,919,676	690,512
8	1,3-dimethyl-2-nitrobe	8:00	480	Auto	1,010,027	506,893
9	triphenyl phosphate (S	22:09	1329	Auto	1,291,058	454,534
10	perylene d-12 (SURR)	27:07	1627	Auto	852,619	145,489
11	hexachlorocyclopentadi	10:11	611	Auto	1,470	1,470
12	bis(2-ethylhexyl)adipa	22:03	1323	Auto	38,399	2,981
13	bis(2-ethylhexyl)phtha	23:02	1382	Auto	347,446	108,008
14	2,6-dinitrotoluene	12:00	720	Auto	76,192	39,126
15	4,4'-dde	20:48	1248	Auto	3,883	2,188

reviewed IS/SURR's

Quantitation Report Quanfile: 25049J Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 11/09/01; SAMPLE 25049
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	740	12:20	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1019	16:59	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1364	22:44	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1253	20:53	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	740	12:20	BB	3.225	UG/L
6	phenanthrene d-10(SURR	A	1019	16:59	BV	3.293	UG/L
7	chrysene d-12 (SURR)	A	1364	22:44	BB	3.480	UG/L
8	1,3-dimethyl-2-nitrobe	A	480	8:00	BV	4.952	UG/L
9	triphenyl phosphate (S	A	1329	22:09	BB	5.106	UG/L
10	perylene d-12 (SURR)	A	1627	27:07	BB	3.273	UG/L
11	hexachlorocyclopentadi	A	611	10:11	BB	0.019	UG/L
16	bis(2-ethylhexyl)adipa	A	1323	22:03	MM	0.186	UG/L
17	bis(2-ethylhexyl)phtha	A	1382	23:02	VB	0.543	UG/L
20	2,6-dinitrotoluene	A	720	12:00	BB	0.711	UG/L
28	4,4'-dde	A	1248	20:48	BB	0.015	UG/L

Comments for Sample AD25049 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-06-2001 Time: 11:11:39

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