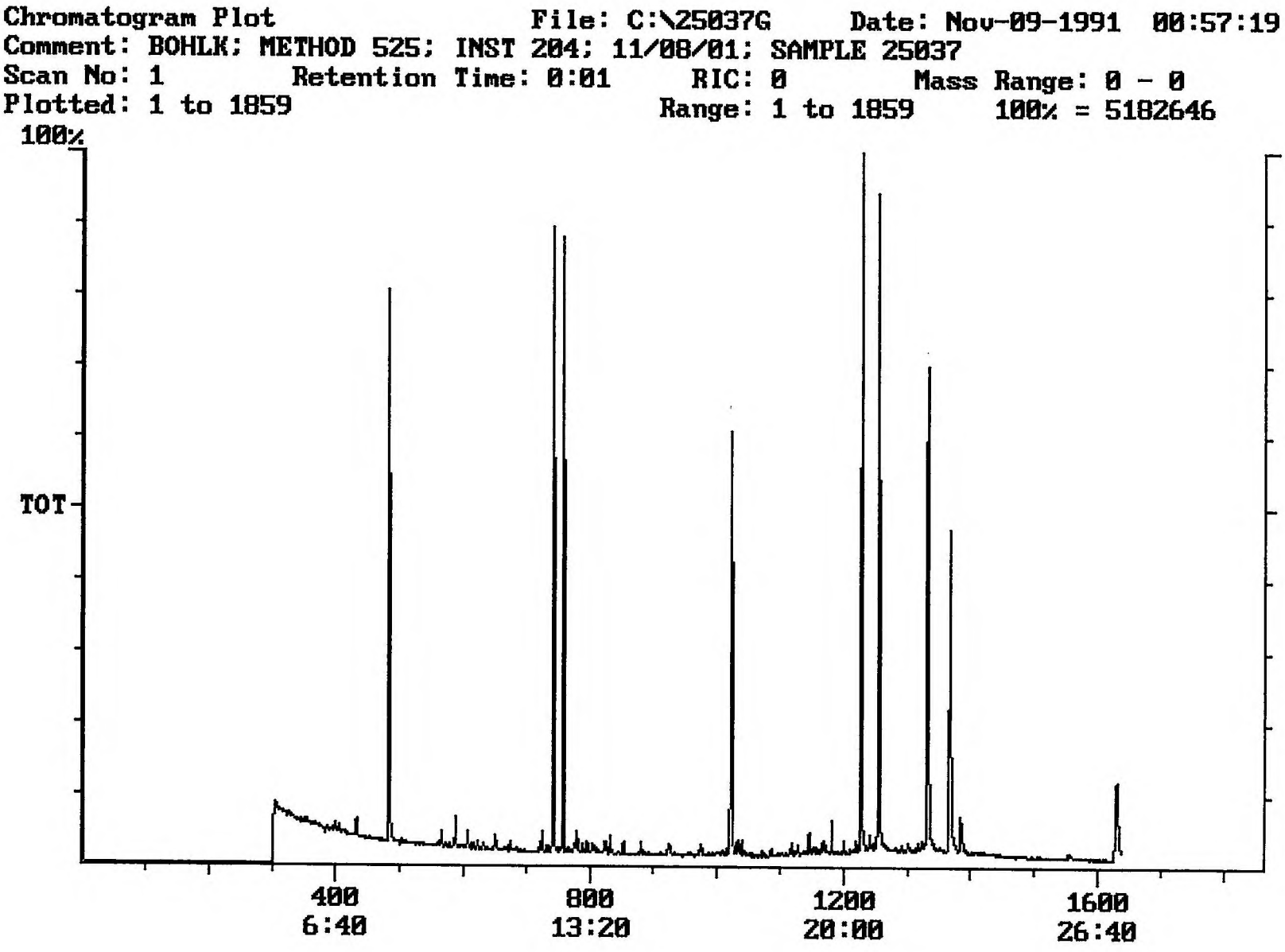
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 11/24/2001 Time: 14:59:43

Sample: AD25037
Analysis: S525 Organic Chemicals - 525.2
Units: ug
Started: 10/24/01 14:58 Ended: 11/8/01 14:58 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		0.10
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		5.1		1.0
20	Surr_Triphenyl phosphate		5.7		1.0
21	Surr_Perylene-d12		3.3		1.0



Integration Report Quanfile: 25037G Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/08/01; SAMPLE 25037
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:22	742	Auto	1,713,541	1,093,465
2	phenanthrene d-10 (IS)	17:02	1022	Auto	2,588,882	993,998
3	chrysene d-12(IS)	22:47	1367	Auto	1,818,602	656,415
4	p-terphenyl d-14(IS)	20:54	1254	Auto	3,298,814	1,488,237
5	acenaphthene d-10(SURR	12:22	742	Auto	1,713,541	1,093,465
6	phenanthrene d-10(SURR	17:02	1022	Auto	2,588,882	993,998
7	chrysene d-12 (SURR)	22:47	1367	Auto	1,818,602	656,415
8	1,3-dimethyl-2-nitrobe	8:02	482	Auto	1,090,369	621,100
9	triphenyl phosphate (S	22:11	1331	Auto	1,353,592	507,315
10	perylene d-12 (SURR)	27:10	1630	Auto	807,661	136,704
11	hexachlorocyclopentadi	10:13	613	Auto	966	966
12	bis(2-ethylhexyl)adipa	22:03	1323	Auto	27,993	1,952
13	bis(2-ethylhexyl)phtha	23:04	1384	Auto	286,533	75,758
14	butachlor	20:29	1229	Auto	0	0

reserved IS/ surr's

Quantitation Report Quanfile: 25037G Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/08/01; SAMPLE 25037
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	742	12:22	BV	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1022	17:02	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1367	22:47	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1254	20:54	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	742	12:22	BV	3.484	UG/L
6	phenanthrene d-10(SURR	A	1022	17:02	BV	3.501	UG/L
7	chrysene d-12 (SURR)	A	1367	22:47	BB	3.416	UG/L
8	1,3-dimethyl-2-nitrobe	A	482	8:02	BV	5.127	UG/L
9	triphenyl phosphate (S	A	1331	22:11	BV	5.651	UG/L
10	perylene d-12 (SURR)	A	1630	27:10	BB	3.273	UG/L
11	hexachlorocyclopentadi	A	613	10:13	BB	0.012	UG/L
16	bis(2-ethylhexyl)adipa	A	1323	22:03	MM	0.143	UG/L
17	bis(2-ethylhexyl)phtha	A	1384	23:04	VB	0.472	UG/L
27	butachlor	A	1229	20:29	BB	0.001	UG/L

Comments for Sample AD25037 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 11-28-2001 Time: 14:29:19

Recoveries for 5 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.