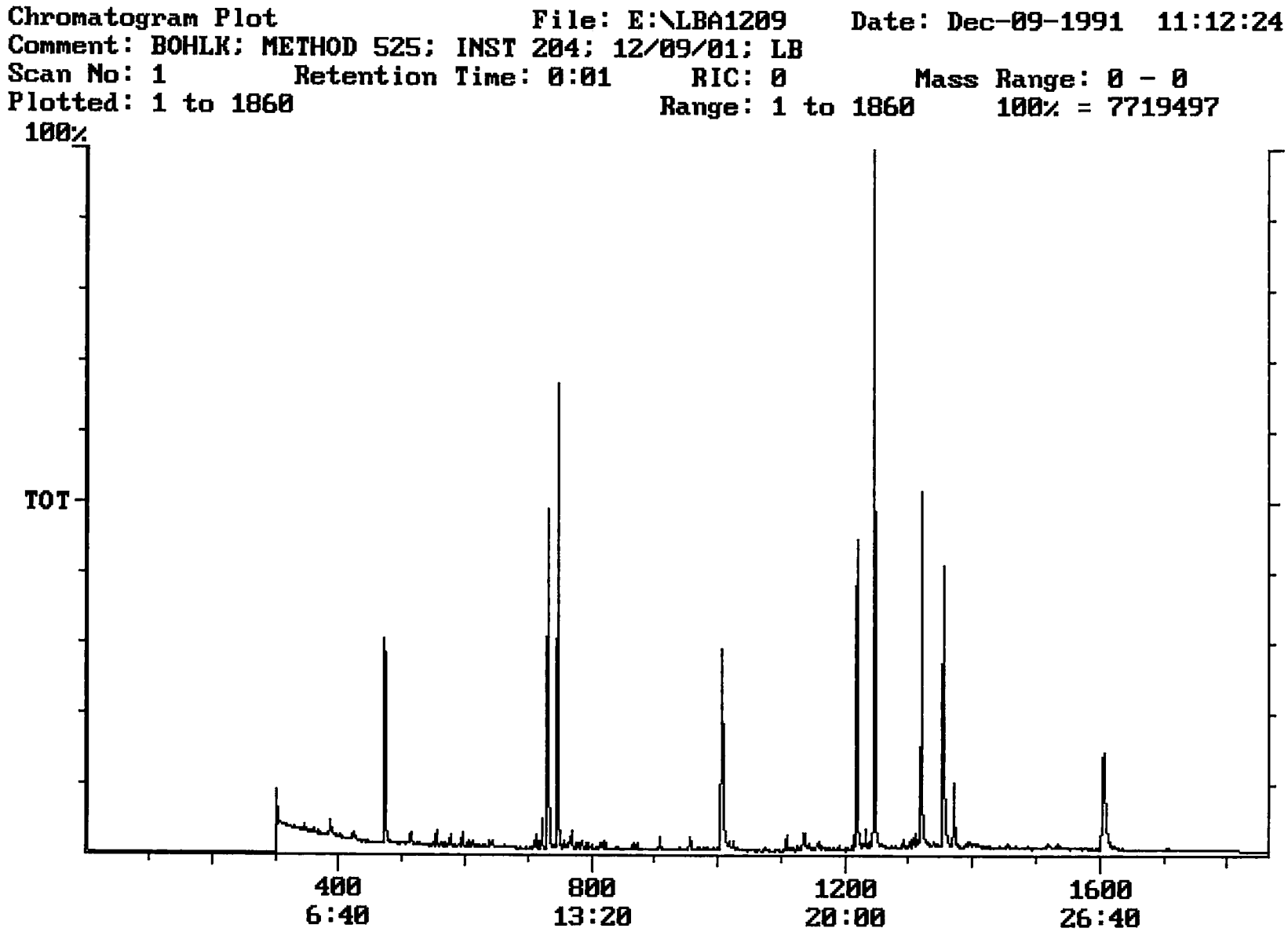
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 01/08/2002 Time: 15:03:06

Sample: AD27769
Analysis: $B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 11/24/01 14:57 Ended: 12/9/01 14:57 Analyst: TB
Result source: manual entry
Page: 1

	Component Name	Vial	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	[illegible]		[illegible]		[illegible]
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		5.8		
20	Surr_Triphenyl phosphate		5.2		
21	Surr_Perylene-d12		4.9		



Integration Report Quanfile: LBA1209 Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 12/09/01; LB
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:11	731	Auto	1,714,932	928,857
2	phenanthrene d-10 (IS)	16:46	1006	Auto	2,633,832	737,756
3	chrysene d-12(IS)	22:34	1354	Auto	2,278,422	977,270
4	p-terphenyl d-14(IS)	20:45	1245	Auto	5,296,038	2,624,894
5	acenaphthene d-10(SURR	12:11	731	Auto	1,714,932	928,857
6	phenanthrene d-10(SURR	16:46	1006	Auto	2,633,832	737,756
7	chrysene d-12 (SURR)	22:34	1354	Auto	2,278,422	977,270
8	1,3-dimethyl-2-nitrobe	7:52	472	Auto	921,922	378,623
9	triphenyl phosphate (S	22:00	1320	Auto	1,148,091	477,298
10	perylene d-12 (SURR)	26:45	1605	Auto	1,614,316	277,899
11	hexachlorobenzene	15:12	912	Auto	5,275	2,436
12	bis(2-ethylhexyl)adipa	21:52	1312	Auto	11,512	3,794
13	bis(2-ethylhexyl)phtha	22:51	1371	Auto	448,060	210,952
14	butachlor	20:18	1218	Auto	12,462	5,838
15	4,4'-dde	20:41	1241	Auto	4,272	4,272

reviewed IS/surr's

Quantitation Report Quanfile: LBA1209 Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 12/09/01; LB
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	731	12:11	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1006	16:46	BB	5.000	UG/L
3	chrysene d-12(IS)	I	1354	22:34	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1245	20:45	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	731	12:11	BB	1.783	UG/L
6	phenanthrene d-10(SURR	A	1006	16:46	BB	2.140	UG/L
7	chrysene d-12 (SURR)	A	1354	22:34	BB	2.262	UG/L
8	1,3-dimethyl-2-nitrobe	A	472	7:52	BV	5.761	UG/L
9	triphenyl phosphate (S	A	1320	22:00	BB	5.200	UG/L
10	perylene d-12 (SURR)	A	1605	26:45	BB	4.863	UG/L
12	hexachlorobenzene	A	912	15:12	BB	0.031	UG/L
16	bis(2-ethylhexyl)adipa	A	1312	21:52	BB	0.034	UG/L
17	bis(2-ethylhexyl)phtha	A	1371	22:51	VV	0.681	UG/L
27	butachlor	A	1218	20:18	BB	0.052	UG/L
28	4,4'-dde	A	1241	20:41	BB	0.019	UG/L