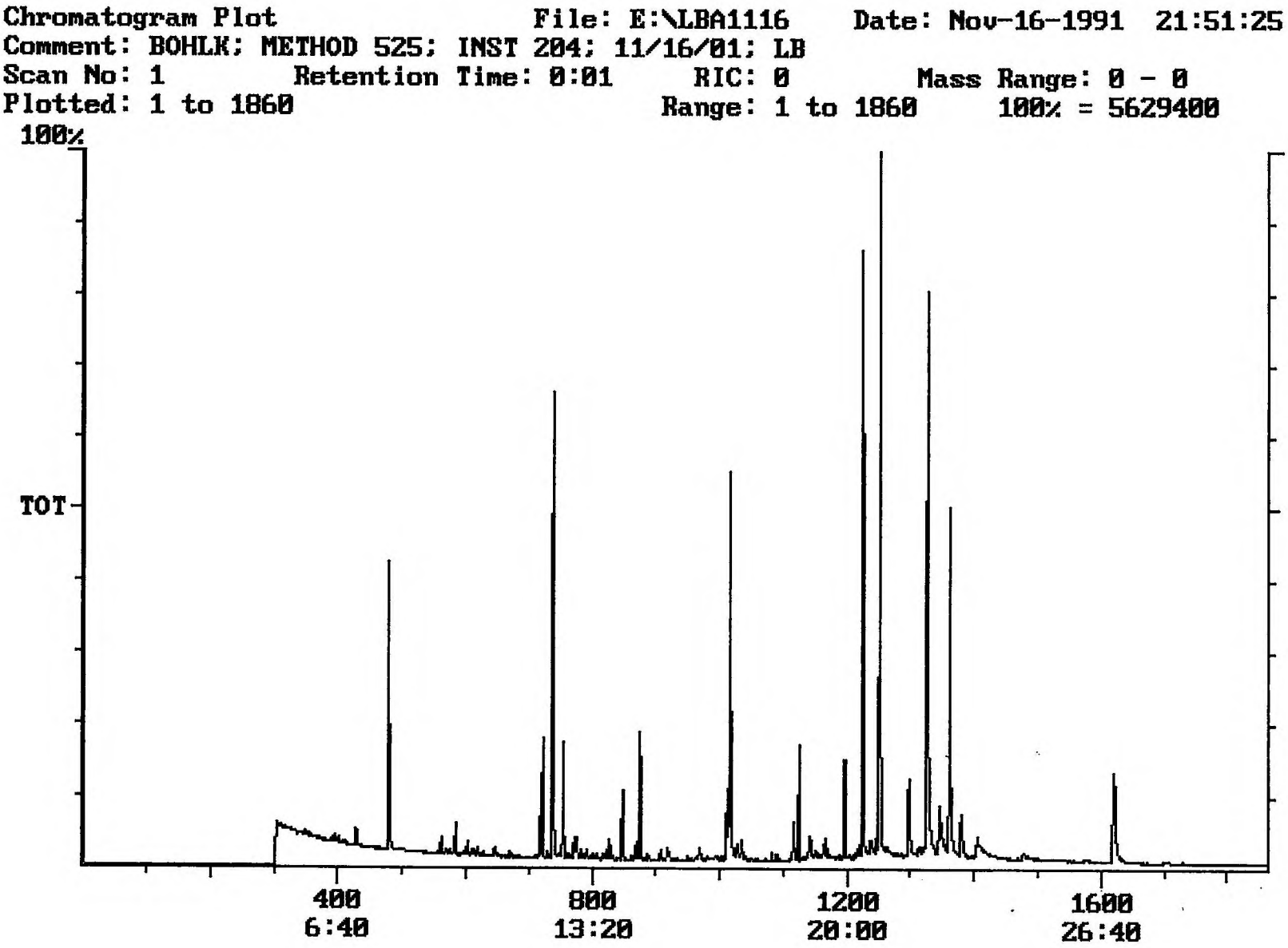
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/08/2001 Time: 17:30:00

Sample: AD25843
Analysis: $B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 11/16/01 13:51 Ended: 11/16/01 13:51 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		4.6		
20	Surr_Triphenyl phosphate		5.3		
21	Surr_Perylene-d12		4.1		



Integration Report Quanfile: LBA1116 Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/16/01; LB
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	1,455,002	827,008
2	phenanthrene d-10 (IS)	16:56	1016	Auto	2,603,126	1,024,944
3	chrysene d-12(IS)	22:41	1361	Auto	2,054,620	836,457
4	p-terphenyl d-14(IS)	20:51	1251	Auto	2,540,657	1,402,124
5	acenaphthene d-10(SURR	12:18	738	Auto	1,455,002	827,008
6	phenanthrene d-10(SURR	16:56	1016	Auto	2,603,126	1,024,944
7	chrysene d-12 (SURR)	22:41	1361	Auto	2,054,620	836,457
8	1,3-dimethyl-2-nitrobe	7:58	478	Auto	835,307	363,699
9	triphenyl phosphate (S	22:06	1326	Auto	1,570,984	674,830
10	perylene d-12 (SURR)	27:00	1620	Auto	1,016,867	166,897
11	hexachlorobenzene	15:21	921	Auto	9,135	3,360
12	bis(2-ethylhexyl)phtha	22:59	1379	Auto	374,962	93,657
13	2,6-dinitrotoluene	11:58	718	Auto	86,221	41,770
14	4,4'-dde	20:46	1246	Auto	11,541	7,414

Reviewed 1S/Sur

Quantitation Report Quanfile: LBA1116 Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/16/01; 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	738	12:18	BB	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	BB	3.768	UG/L
6	phenanthrene d-10(SURR	A	1016	16:56	BV	4.397	UG/L
7	chrysene d-12 (SURR)	A	1361	22:41	BB	5.203	UG/L
8	1,3-dimethyl-2-nitrobe	A	478	7:58	BV	4.609	UG/L
9	triphenyl phosphate (S	A	1326	22:06	BB	5.334	UG/L
10	perylene d-12 (SURR)	A	1620	27:00	BB	4.101	UG/L
12	hexachlorobenzene	A	921	15:21	BB	0.046	UG/L
17	bis(2-ethylhexyl)phtha	A	1379	22:59	VB	0.467	UG/L
20	2,6-dinitrotoluene	A	718	11:58	BB	0.927	UG/L
28	4,4'-dde	A	1246	20:46	BB	0.052	UG/L