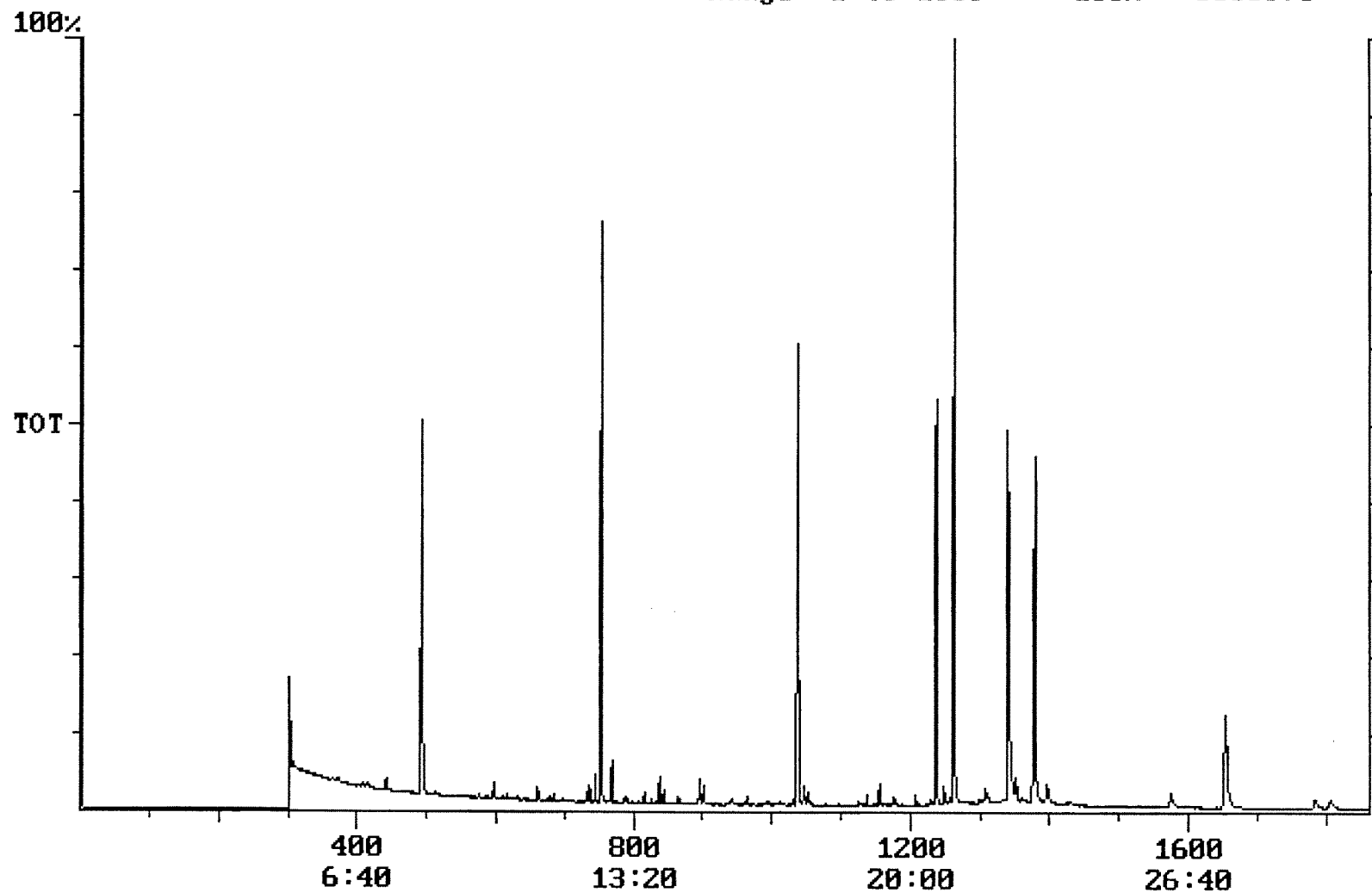


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/05/2001 Time: 08:55:08

Sample: AD22754
Analysis: \$B1525 Organic Chemicals - 525.2; lab blank
Units: ug
Started: 9/25/01 08:50 Ended: 10/1/01 08:50 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		0.07
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-nitrobenzen		4.6		
20	Surr_Triphenyl phosphate		5.0		
21	Surr_Perylene-d12		4.3		

Chromatogram Plot File: E:\LBA1001 Date: Oct-01-1991 20:05:50
Comment: BOHLK; METHOD 525; INST 204; 10/01/01; LB
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1859 Range: 1 to 1859 100% = 5635970



Integration Report

Quanfile: LBA1001

Quan Entries: 15

Comment: BOHLK; METHOD 525; INST 204; 10/01/01; LB

Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10 (IS)	12:33	753	Auto	1,480,751	1,018,627
2	phenanthrene d-10 (IS)	17:17	1037	Auto	2,256,944	1,148,170
3	chrysene d-12 (IS)	22:58	1378	Auto	1,639,313	724,560
4	p-terphenyl d-14 (IS)	21:02	1262	Auto	2,535,475	1,809,415
5	acenaphthene d-10 (SURR)	12:33	753	Auto	1,480,751	1,018,627
6	phenanthrene d-10 (SURR)	17:17	1037	Auto	2,256,944	1,148,170
7	chrysene d-12 (SURR)	22:58	1378	Auto	1,639,313	724,560
8	1,3-dimethyl-2-nitrobenzene	8:13	493	Auto	896,201	425,912
9	triphenyl phosphate (S)	22:20	1340	Auto	1,110,888	421,363
10	perylene d-12 (SURR)	27:33	1653	Auto	870,927	165,115
11	bis(2-ethylhexyl)adipate	22:20	1340	Auto	14,931	3,225
12	bis(2-ethylhexyl)phthalate	23:15	1395	Auto	202,448	39,831
13	2,6-dinitrotoluene	12:12	732	Auto	17,932	10,781
14	2,4-dinitrotoluene	13:19	799	Auto	1,086	1,086
15	4,4'-dinitrodiphenyl ether	20:58	1258	Auto	1,576	1,576

Quantitation Report

Quanfile: LBA1001

Quan Entries: 15

Comment: BOHLK; METHOD 525; INST 204; 10/01/01; 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	753	12:33	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1037	17:17	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1378	22:58	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1262	21:02	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	753	12:33	BB	3.901	UG/L
6	phenanthrene d-10(SURR	A	1037	17:17	BV	3.877	UG/L
7	chrysene d-12 (SURR)	A	1378	22:58	BB	4.154	UG/L
8	1,3-dimethyl-2-nitrobe	A	493	8:13	BB	4.641	UG/L
9	triphenyl phosphate (S	A	1340	22:20	BB	5.033	UG/L
10	perylene d-12 (SURR)	A	1653	27:33	BB	4.277	UG/L
16	bis(2-ethylhexyl)adipa	A	1340	22:20	MM	0.056	UG/L
17	bis(2-ethylhexyl)phtha	A	1395	23:15	MM	0.431	UG/L
20	2,6-dinitrotoluene	A	732	12:12	BB	0.175	UG/L
21	2,4-dinitrotoluene	A	799	13:19	BB	0.009	UG/L
28	4,4'-dde	A	1258	20:58	BB	0.008	UG/L
<i>reviewed IS/surrs</i>							

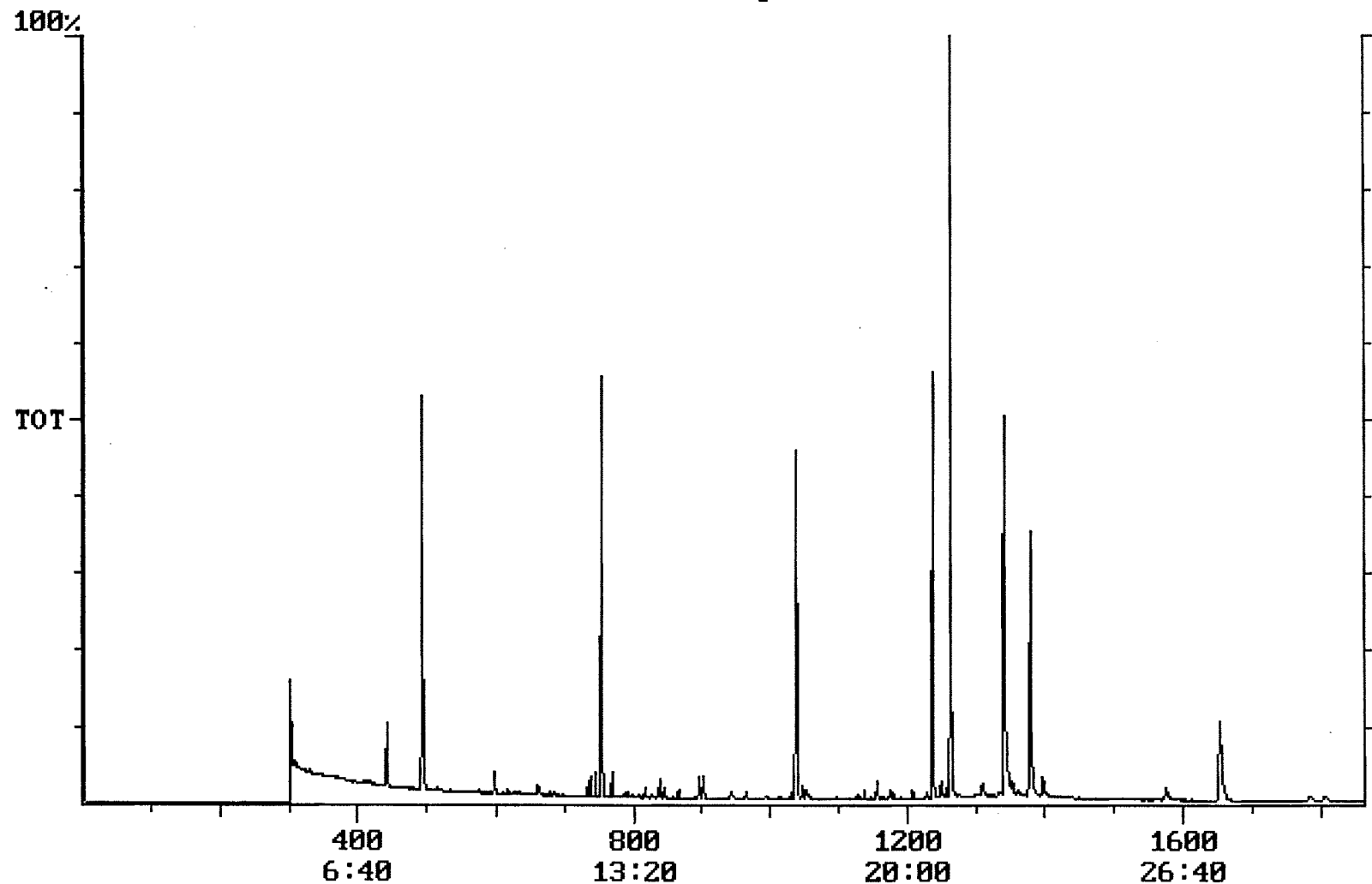
Chromatogram Plot

File: E:\LBA1002 Date: Oct-02-1991 19:48:53

Comment: BOHLK; METHOD 525; INST 204; 10/02/01; LB REINJD

Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0

Plotted: 1 to 1859 Range: 1 to 1859 100% = 6787045



Integration Report

Quanfile: LBA1002

Quan Entries: 14

Comment: BOHLK; METHOD 525; INST 204; 10/02/01; LB REINJD

Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:34	754	Auto	1,597,356	862,500
2	phenanthrene d-10 (IS)	17:17	1037	Auto	2,461,852	1,055,514
3	chrysene d-12(IS)	22:59	1379	Auto	1,799,925	653,308
4	p-terphenyl d-14(IS)	21:03	1263	Auto	2,931,403	2,139,062
5	acenaphthene d-10(SURR)	12:34	754	Auto	1,597,356	862,500
6	phenanthrene d-10(SURR)	17:17	1037	Auto	2,461,852	1,055,514
7	chrysene d-12 (SURR)	22:59	1379	Auto	1,799,925	653,308
8	1,3-dimethyl-2-nitrobenzophenone	8:13	493	Auto	984,766	550,231
9	triphenyl phosphate (S)	22:21	1341	Auto	1,236,907	503,133
10	perylene d-12 (SURR)	27:33	1653	Auto	962,026	182,878
11	bis(2-ethylhexyl)adipate	22:21	1341	Auto	21,792	4,084
12	bis(2-ethylhexyl)phthalate	23:16	1396	Auto	237,619	50,255
13	2,6-dinitrotoluene	12:12	732	Auto	24,597	11,789
14	4,4'-dichlorodiphenyl ether	20:59	1259	Auto	2,379	1,378

Quantitation Report

Quanfile: LBA1002

Quan Entries: 14

Comment: BOHLK; METHOD 525; INST 204; 10/02/01; LB REINJD

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	754	12:34	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1037	17:17	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1379	22:59	BV	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1263	21:03	BB	5.000	UG/L
5	acenaphthene d-10(SURR)	A	754	12:34	BB	3.640	UG/L
6	phenanthrene d-10(SURR)	A	1037	17:17	BV	3.658	UG/L
7	chrysene d-12 (SURR)	A	1379	22:59	BV	3.945	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	493	8:13	BB	4.728	UG/L
9	triphenyl phosphate (S)	A	1341	22:21	BB	5.104	UG/L
10	perylene d-12 (SURR)	A	1653	27:33	BB	4.302	UG/L
16	bis(2-ethylhexyl)adipate	A	1341	22:21	MM	0.075	UG/L
17	bis(2-ethylhexyl)phthalate	A	1396	23:16	VB	0.461	UG/L
20	2,6-dinitrotoluene	A	732	12:12	BB	0.222	UG/L
28	4,4'-dichlorodiphenyl ether	A	1259	20:59	BB	0.011	UG/L
<i>reversed IS/surrs</i>							