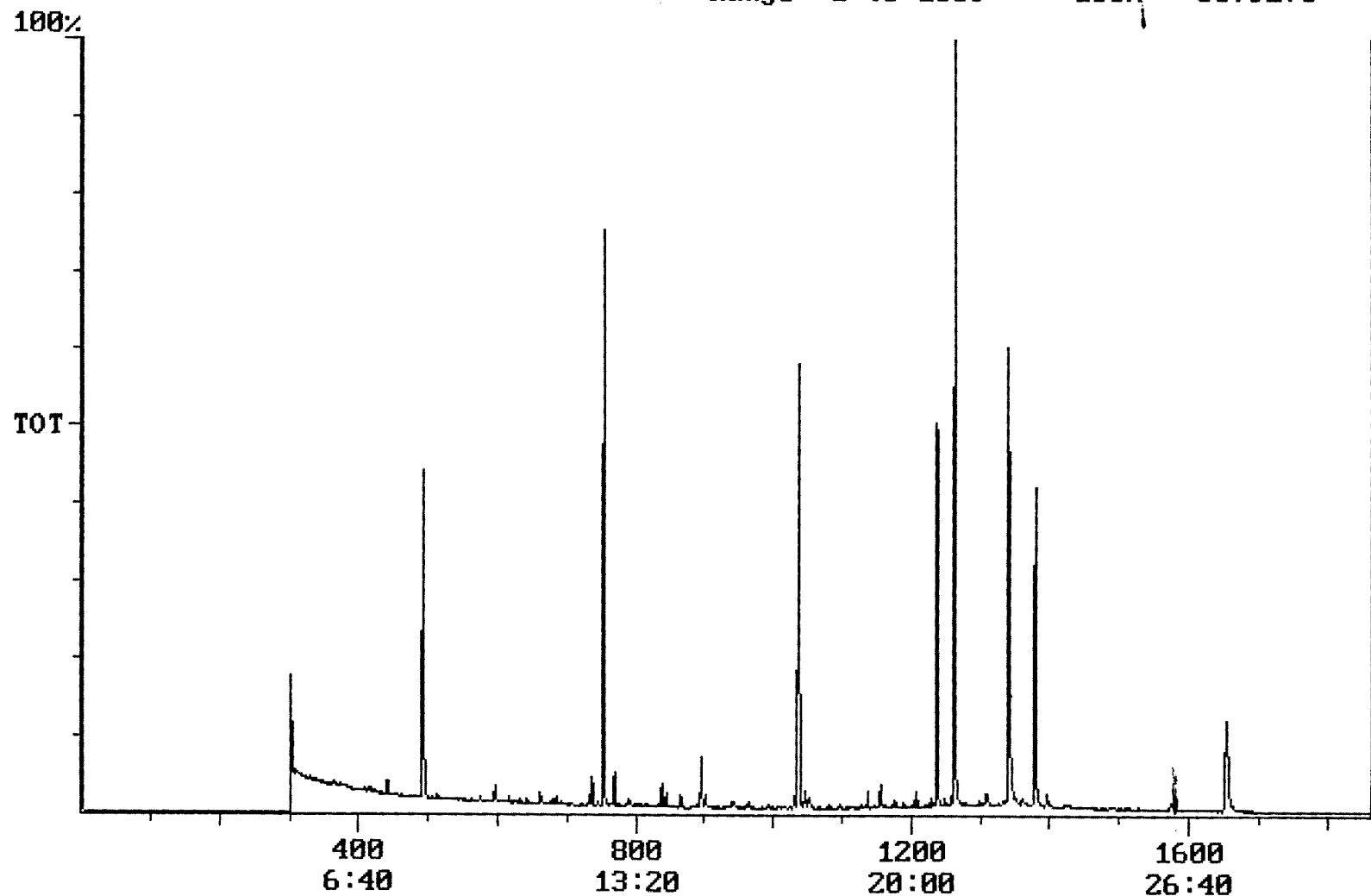


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/05/2001 Time: 10:16:59

Sample: AD22754
Analysis: \$525 Organic Chemicals - 525.2
Units: ug
Started: 9/25/01 10:15 Ended: 10/1/01 10:15 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		1.0
20	Surr_Triphenyl phosphate		5.9		1.0
21	Surr_Perylene-d12		4.2		1.0

Chromatogram Plot File: E:\22754AAA Date: Oct-01-1991 21:14:41
 Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22754
 Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
 Plotted: 1 to 1859 Range: 1 to 1859 100% = 5879278



Integration Report Quanfile: 22754AAA Quan Entries: 17
 Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22754
 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	1,507,865	1,065,721
2	phenanthrene d-10 (IS)	17:17	1037	Auto	2,389,131	1,148,451
3	chrysene d-12(IS)	22:58	1378	Auto	1,656,795	702,439
4	p-terphenyl d-14(IS)	21:02	1262	Auto	2,575,946	1,847,666
5	acenaphthene d-10(SURR)	12:33	753	Auto	1,507,865	1,065,721
6	phenanthrene d-10(SURR)	17:17	1037	Auto	2,389,131	1,148,451
7	chrysene d-12 (SURR)	22:58	1378	Auto	1,656,795	702,439
8	1,3-dimethyl-2-nitrobe	8:12	492	Auto	908,221	380,063
9	triphenyl phosphate (S	22:20	1340	Auto	1,312,201	541,858
10	perylene d-12 (SURR)	27:33	1653	Auto	856,828	169,894
11	hexachlorocyclopentadi	10:21	621	Auto	0	0
12	hexachlorobenzene	15:44	944	Auto	5,242	2,245
13	bis(2-ethylhexyl)adipa	22:20	1340	Auto	16,037	5,916
14	bis(2-ethylhexyl)phtha	23:15	1395	Auto	115,191	26,657
15	2,6-dinitrotoluene	12:12	732	Auto	20,452	12,617
16	2,4-dinitrotoluene	13:19	799	Auto	1,413	1,413
17	4,4'-dde	20:58	1258	Auto	5,163	3,658

Quantitation Report Quanfile: 22754AAA Quan Entries: 17
 Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22754
 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	3.910	UG/L
6	phenanthrene d-10(SURR)	A	1037	17:17	BV	4.040	UG/L
7	chrysene d-12 (SURR)	A	1378	22:58	BB	4.132	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	492	8:12	BB	4.619	UG/L
9	triphenyl phosphate (S)	A	1340	22:20	BB	5.882	UG/L
10	perylene d-12 (SURR)	A	1653	27:33	BB	4.163	UG/L
11	hexachlorocyclopentadiene	A	621	10:21	BB	0.001	UG/L
12	hexachlorobenzene	A	944	15:44	BB	0.032	UG/L
16	bis(2-ethylhexyl)adipate	A	1340	22:20	MM	0.060	UG/L
17	bis(2-ethylhexyl)phthalate	A	1395	23:15	MM	0.244	UG/L
20	2,6-dinitrotoluene	A	732	12:12	BB	0.195	UG/L
21	2,4-dinitrotoluene	A	799	13:19	BB	0.011	UG/L
28	4,4'-dichlorodiphenyl ether	A	1258	20:58	BB	0.024	UG/L
<i>revised IS/surrogate</i>							

Comments for Sample AD22754 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 10-08-2001 Time: 12:51:27

Recoveries for three of the eighteen compounds in the MDL Standard were outside of their acceptance ranges. Recoveries for six of the eighteen compounds in the LCS Standard were outside of their acceptance ranges.