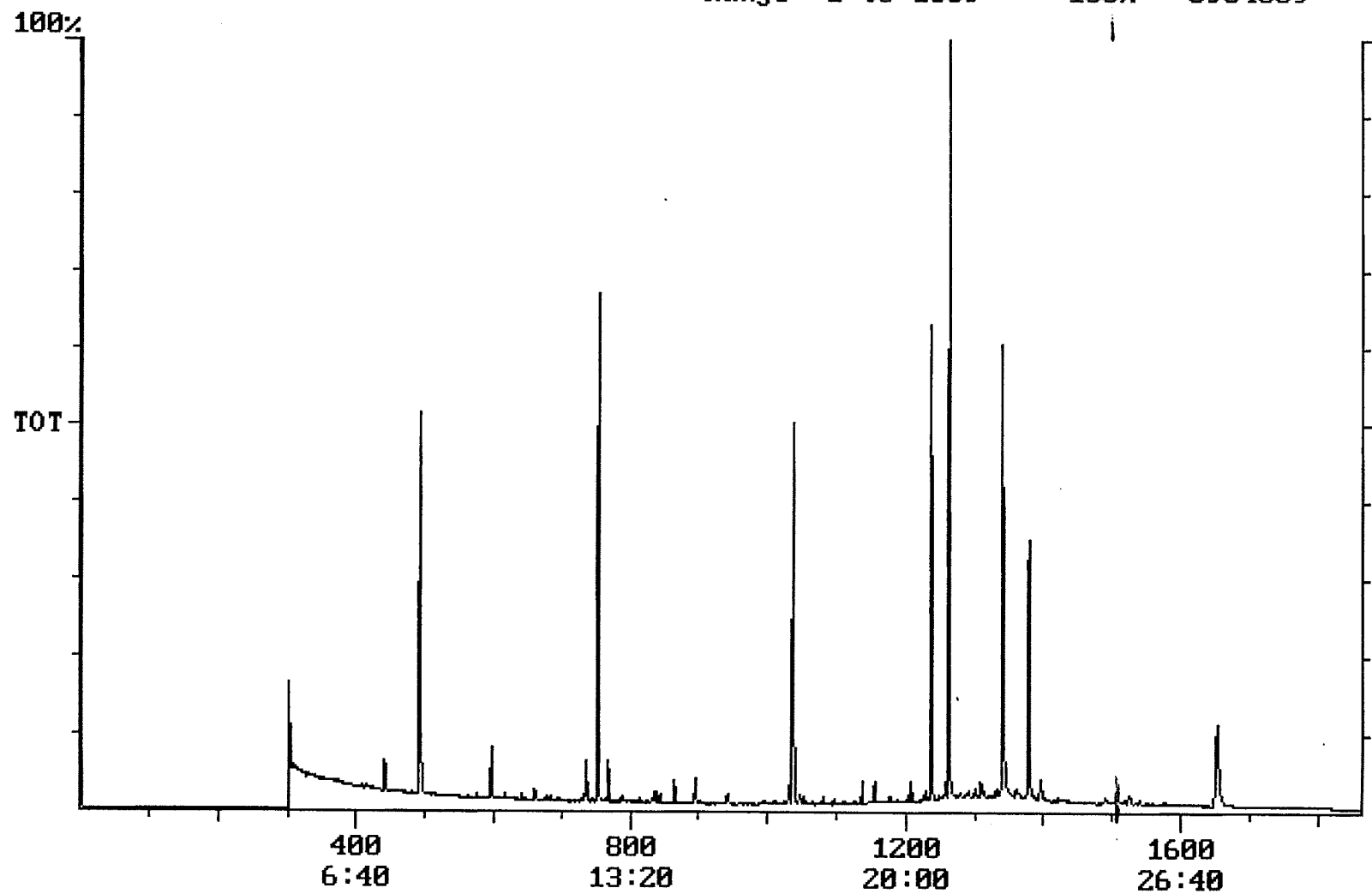


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/04/2001 Time: 10:16:52

Sample: AD22752
Analysis: \$525 Organic Chemicals - 525.2
Units: ug
Started: 9/25/01 10:01 Ended: 9/30/01 10:01 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		4.9		1.0
20	Surr_Triphenyl phosphate		6.0		1.0
21	Surr_Perylene-d12		4.5		1.0

Chromatogram Plot File: E:\22752CCC Date: Oct-01-1991 03:42:22
Comment: BOHLK; METHOD 525; INST 204; 09/30/01; SAMPLE 22752
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1859 Range: 1 to 1859 100% = 5934339



Integration Report Quanfile: 22752CCC Quan Entries: 13
 Comment: BOHLK; METHOD 525; INST 204; 09/30/01; SAMPLE 22752
 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	1,470,446	919,772
2	phenanthrene d-10 (IS)	17:17	1037	Auto	2,215,051	966,176
3	chrysene d-12 (IS)	22:58	1378	Auto	1,460,527	553,536
4	p-terphenyl d-14 (IS)	21:02	1262	Auto	2,455,563	1,800,324
5	acenaphthene d-10 (SURR)	12:33	753	Auto	1,470,446	919,772
6	phenanthrene d-10 (SURR)	17:17	1037	Auto	2,215,051	966,176
7	chrysene d-12 (SURR)	22:58	1378	Auto	1,460,527	553,536
8	1,3-dimethyl-2-nitrobenzene	8:12	492	Auto	934,280	439,962
9	triphenyl phosphate (S)	22:20	1340	Auto	1,175,423	521,105
10	perylene d-12 (SURR)	27:32	1652	Auto	813,633	163,600
11	bis(2-ethylhexyl)phthalate	23:15	1395	Auto	171,054	47,223
12	2,6-dinitrotoluene	12:12	732	Auto	18,408	9,299
13	2,4-dinitrotoluene	13:19	799	Auto	2,822	1,841

Quan Entries: 13

22752

(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	4.000	UG/L
6	phenanthrene d-10(SURR)	A	1037	17:17	BV	3.929	UG/L
7	chrysene d-12 (SURR)	A	1378	22:58	BB	3.821	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	492	8:12	BB	4.873	UG/L
9	triphenyl phosphate (S)	A	1340	22:20	BB	5.977	UG/L
10	perylene d-12 (SURR)	A	1652	27:32	BB	4.484	UG/L
17	bis(2-ethylhexyl)phthalate	A	1395	23:15	BB	0.409	UG/L
20	2,6-dinitrotoluene	A	732	12:12	BB	0.180	UG/L
21	2,4-dinitrotoluene	A	799	13:19	BB	0.022	UG/L

reversed IS/SURR's

Comments for Sample AD22752 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 10-05-2001 Time: 10:28:26

The recovery for bis(2-Ethylhexyl)phthalate in the MDL Standard was above its acceptance range. The recoveries for Terbacil in the LCS Standard and the Spiked Sample were above their acceptance ranges.