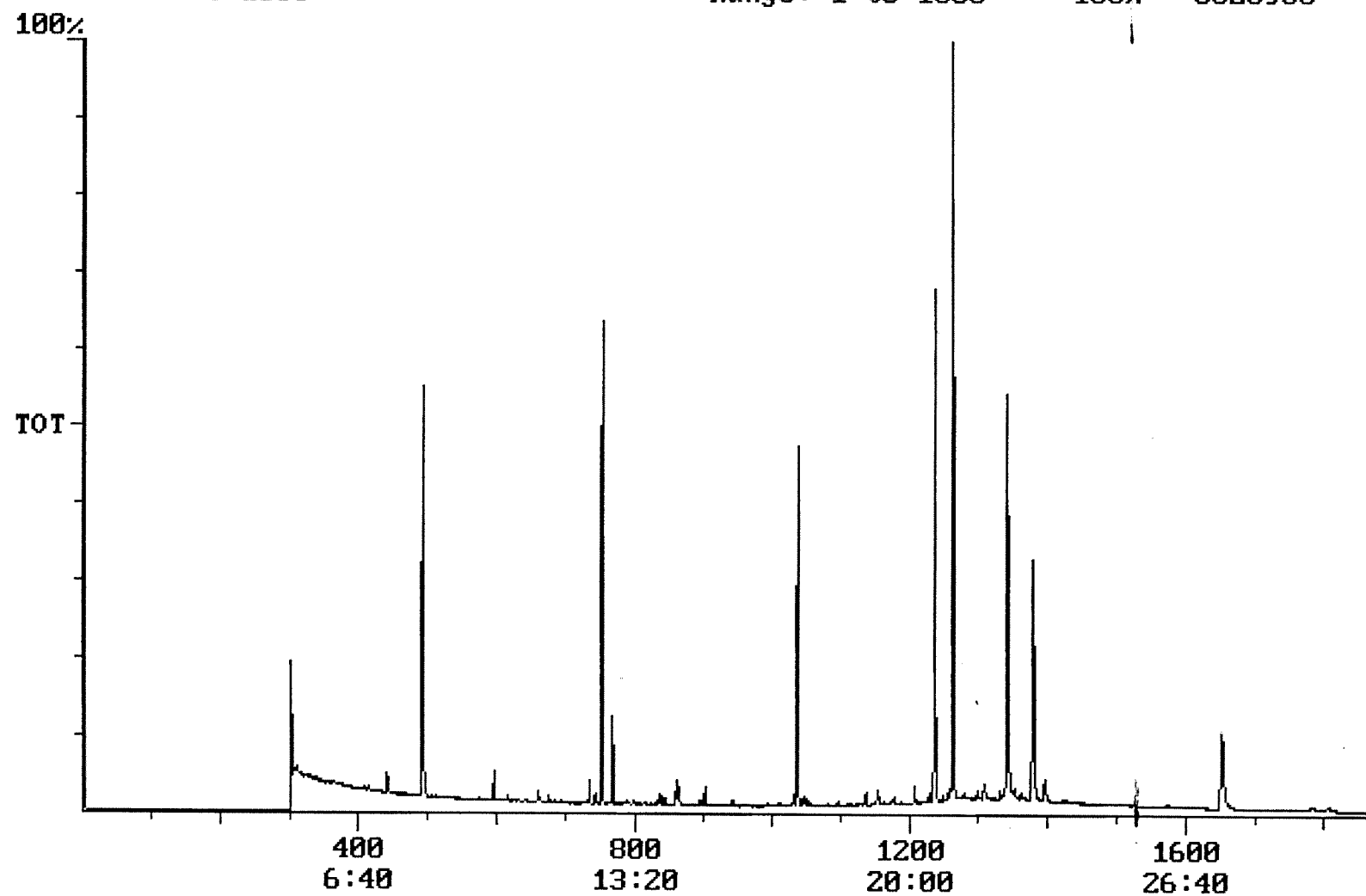


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/04/2001 Time: 10:17:33

Sample: AD22753
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		4.6		1.0
20	Surr_Triphenyl phosphate		6.0		1.0
21	Surr_Perylene-d12		4.3		1.0

Chromatogram Plot File: E:\22753DDD Date: Oct-01-1991 04:16:49
Comment: BOHLK; METHOD 525; INST 204; 09/30/01; SAMPLE 22753
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 6028950



Integration Report

Quanfile: 22753DDD

Quan Entries: 12

Comment: BOHLK; METHOD 525; INST 204; 09/30/01; SAMPLE 22753

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	1,476,721	884,913
2	phenanthrene d-10 (IS)	17:16	1036	Auto	2,207,916	966,826
3	chrysene d-12 (IS)	22:57	1377	Auto	1,464,949	540,314
4	p-terphenyl d-14 (IS)	21:02	1262	Auto	2,565,610	1,870,666
5	acenaphthene d-10 (SURR)	12:33	753	Auto	1,476,721	884,913
6	phenanthrene d-10 (SURR)	17:16	1036	Auto	2,207,916	966,826
7	chrysene d-12 (SURR)	22:57	1377	Auto	1,464,949	540,314
8	1,3-dimethyl-2-nitrobenzene	8:12	492	Auto	883,476	492,137
9	triphenyl phosphate (S)	22:20	1340	Auto	1,183,705	484,215
10	perylene d-12 (SURR)	27:32	1652	Auto	776,927	148,322
11	bis(2-ethylhexyl)phthalate	23:15	1395	Auto	177,762	47,236
12	2,6-dinitrotoluene	12:11	731	Auto	10,080	5,489

Quan Entries: 12

22753

(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	1036	17:16	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1377	22:57	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1262	21:02	BV	5.000	UG/L
5	acenaphthene d-10(SURR)	A	753	12:33	BB	3.844	UG/L
6	phenanthrene d-10(SURR)	A	1036	17:16	BV	3.748	UG/L
7	chrysene d-12 (SURR)	A	1377	22:57	BB	3.669	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	492	8:12	BB	4.588	UG/L
9	triphenyl phosphate (S)	A	1340	22:20	BV	6.001	UG/L
10	perylene d-12 (SURR)	A	1652	27:32	BB	4.269	UG/L
17	bis(2-ethylhexyl)phthalate	A	1395	23:15	MM	0.424	UG/L
20	2,6-dinitrotoluene	A	731	12:11	BB	0.099	UG/L

reviewed IS/Suror's.

Comments for Sample AD22753 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 10-05-2001 Time: 10:37:48

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.