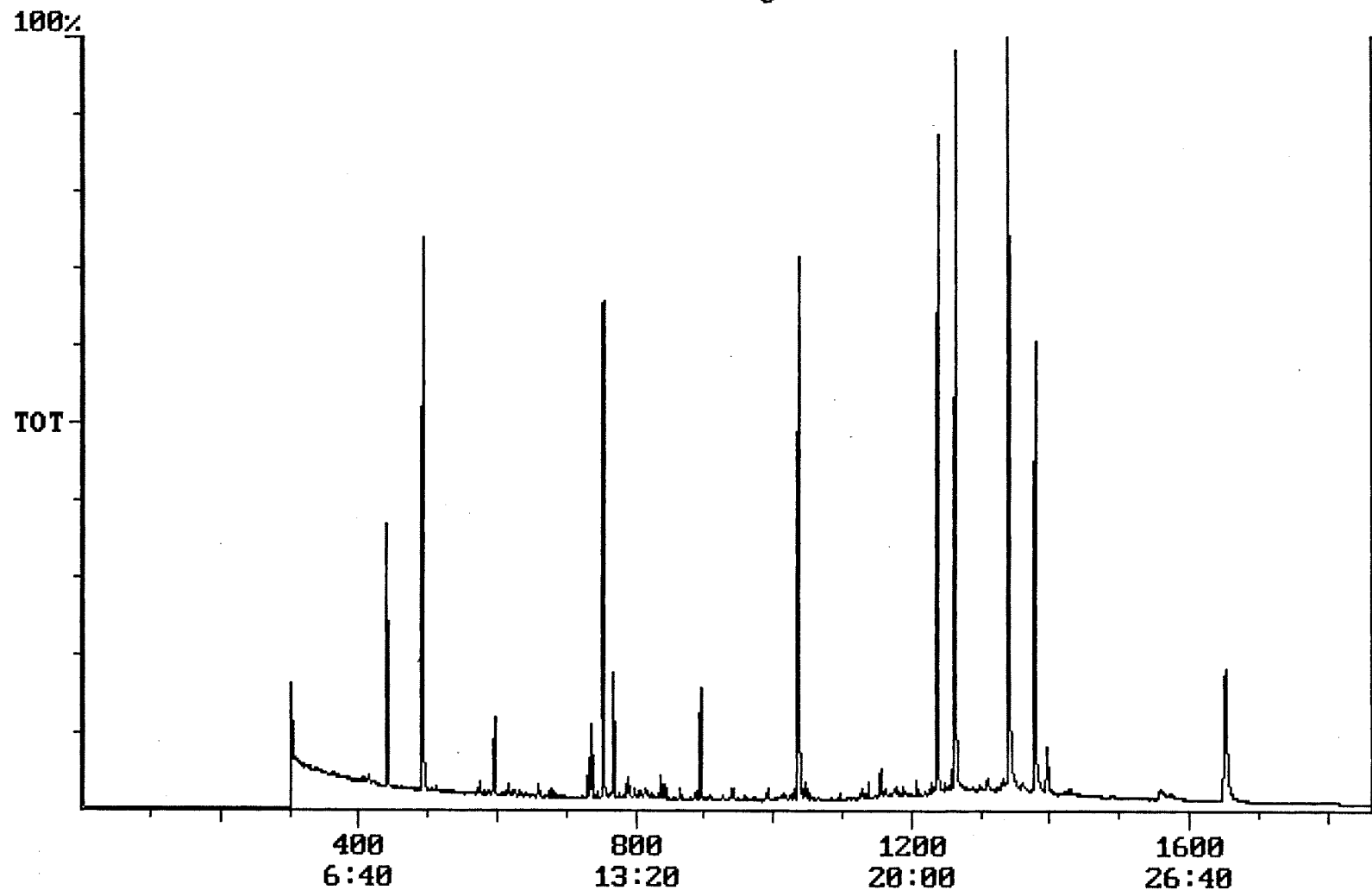


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/24/2001 Time: 11:08:02

Sample: AD23652
Analysis: \$525 Organic Chemicals - 525.2
Units: ug
Started: 10/5/01 11:01 Ended: 10/16/01 11:01 Analyst: TB
Result source: manual entry
Page: 1

	Component Name	Vol	Result	Qualifier	MDL
1	Hexachlorocyclopentadiene		< MDL		0.10
2	Hexachlorobenzene		< MDL		0.10
3	Simazine		< MDL		0.40
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-nitrobenze		5.0		1.0
20	Surr_Triphenyl phosphate		5.7		1.0
21	Surr_Perylene-d12		5.9		1.0

Chromatogram Plot File: E:\23652TTE Date: Oct-17-1991 00:46:45
Comment: BOHLK; METHOD 525; INST 204; 10/16/01; SAMPLE 23652
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1859 Range: 1 to 1859 100% = 6282156



Integration Report

Quanfile: 23652TTE

Quan Entries: 17

Comment: BOHLK; METHOD 525; INST 204; 10/16/01; SAMPLE 23652

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	2,139,176	956,862
2	phenanthrene d-10 (IS)	17:16	1036	Auto	3,292,105	1,528,060
3	chrysene d-12 (IS)	22:58	1378	Auto	2,437,189	1,058,816
4	p-terphenyl d-14 (IS)	21:02	1262	Auto	3,606,462	1,954,205
5	acenaphthene d-10 (SURR)	12:33	753	Auto	2,139,176	956,862
6	phenanthrene d-10 (SURR)	17:16	1036	Auto	3,292,105	1,528,060
7	chrysene d-12 (SURR)	22:58	1378	Auto	2,437,189	1,058,816
8	1,3-dimethyl-2-nitrobenzene	8:12	492	Auto	1,290,852	705,919
9	triphenyl phosphate (S)	22:20	1340	Auto	1,988,764	943,502
10	perylene d-12 (SURR)	27:32	1652	Auto	1,786,898	327,513
11	hexachlorocyclopentadiene	10:23	623	Auto	2,736	1,877
12	hexachlorobenzene	15:43	943	Auto	9,285	4,268
13	bis(2-ethylhexyl)adipate	22:20	1340	Auto	26,532	7,888
14	bis(2-ethylhexyl)phthalate	23:15	1395	Auto	389,539	110,088
15	2,6-dinitrotoluene	12:11	731	Auto	83,146	44,776
16	2,4-dinitrotoluene	13:19	799	Auto	3,667	1,684
17	4,4'-dichlorodiphenyl ether	20:58	1258	Auto	9,137	6,087

Quantitation Report

Quanfile: 23652TTE

Quan Entries: 17

Comment: BOHLK; METHOD 525; INST 204; 10/16/01; SAMPLE 23652

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	1036	17:16	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	BV	5.000	UG/L
5	acenaphthene d-10(SURR	A	753	12:33	BB	4.258	UG/L
6	phenanthrene d-10(SURR	A	1036	17:16	BV	4.448	UG/L
7	chrysene d-12 (SURR)	A	1378	22:58	BB	4.759	UG/L
8	1,3-dimethyl-2-nitrobe	A	492	8:12	BV	5.036	UG/L
9	triphenyl phosphate (S	A	1340	22:20	BV	5.711	UG/L
10	perylene d-12 (SURR)	A	1652	27:32	BB	5.907	UG/L
11	hexachlorocyclopentadi	A	623	10:23	BB	0.045	UG/L
12	hexachlorobenzene	A	943	15:43	BB	0.035	UG/L
16	bis(2-ethylhexyl)adipa	A	1340	22:20	MM	0.066	UG/L
17	bis(2-ethylhexyl)phtha	A	1395	23:15	VB	0.518	UG/L
20	2,6-dinitrotoluene	A	731	12:11	BV	0.550	UG/L
21	2,4-dinitrotoluene	A	799	13:19	BB	0.021	UG/L
28	4,4'-dde	A	1258	20:58	BB	0.033	UG/L

reversed IS/standards

Comments for Sample AD23652 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 11-04-2001 Time: 12:04:11

The recovery for Terbacil in the LCS Standard was above its acceptance range. Recoveries for four and three of the eighteen compounds in the Spiked Sample and the Spiked Duplicate Sample respectively, were outside of their acceptance ranges.