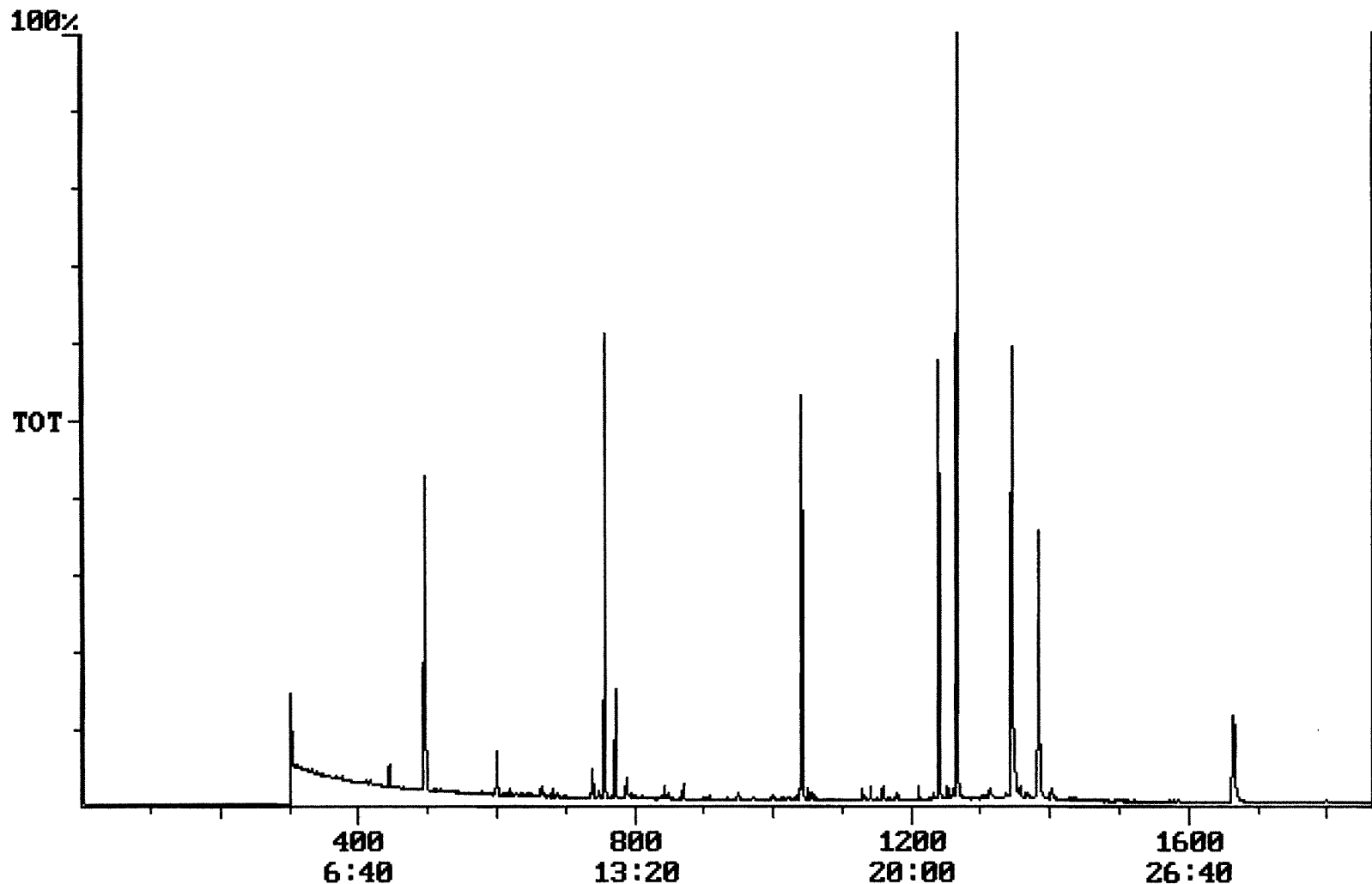


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 09/25/2001 Time: 17:24:04

Sample: AD22495
Analysis: \$525 Organic Chemicals - 525.2
Units: ug
Started: 9/19/01 17:14 Ended: 9/21/01 17:14 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		5.1		1.0
20	Surr_Triphenyl phosphate		6.0		1.0
21	Surr_Perylene-d12		4.7		1.0

Chromatogram Plot File: E:\22495B Date: Sep-21-1991 17:08:45
Comment: BOHLK; METHOD 525; INST 204; 09/21/01; SAMPLE 22495
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 3778281



Integration Report Quanfile: 22495B Quan Entries: 13
 Comment: BOHLK; METHOD 525; INST 204; 09/21/01; SAMPLE 22495
 Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:37	757	Auto	832,134	541,166
2	phenanthrene d-10 (IS)	17:22	1042	Auto	1,353,040	677,215
3	chrysene d-12(IS)	23:03	1383	Auto	988,191	379,729
4	p-terphenyl d-14(IS)	21:06	1266	Auto	1,725,226	1,268,041
5	acenaphthene d-10(SURR)	12:37	757	Auto	832,134	541,166
6	phenanthrene d-10(SURR)	17:22	1042	Auto	1,353,040	677,215
7	chrysene d-12 (SURR)	23:03	1383	Auto	988,191	379,729
8	1,3-dimethyl-2-nitrobenzene	8:16	496	Auto	521,833	242,050
9	triphenyl phosphate (S)	22:25	1345	Auto	825,348	341,664
10	perylene d-12 (SURR)	27:44	1664	Auto	595,177	111,286
11	bis(2-ethylhexyl)phthalate	23:21	1401	Auto	77,855	15,301
12	2,6-dinitrotoluene	12:15	735	Auto	6,484	4,189
13	2,4-dinitrotoluene	13:23	803	Auto	1,137	1,137

Quantitation Report Quanfile: 22495B Quan Entries: 13
 Comment: BOHLK; METHOD 525; INST 204; 09/21/01; SAMPLE 22495
 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	757	12:37	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1042	17:22	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1383	23:03	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1266	21:06	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	757	12:37	BB	3.337	UG/L
6	phenanthrene d-10(SURR	A	1042	17:22	BV	3.654	UG/L
7	chrysene d-12 (SURR)	A	1383	23:03	BB	3.729	UG/L
8	1,3-dimethyl-2-nitrobe	A	496	8:16	BB	5.120	UG/L
9	triphenyl phosphate (S	A	1345	22:25	BV	5.969	UG/L
10	perylene d-12 (SURR)	A	1664	27:44	BB	4.693	UG/L
17	bis(2-ethylhexyl)phtha	A	1401	23:21	MM	0.348	UG/L
20	2,6-dinitrotoluene	A	735	12:15	BB	0.127	UG/L
21	2,4-dinitrotoluene	A	803	13:23	BB	0.018	UG/L

Comments for Sample AD22495 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 09-26-2001 Time: 15:32:30

The recovery for Terbacil in the LCS Standard was above its acceptance range.