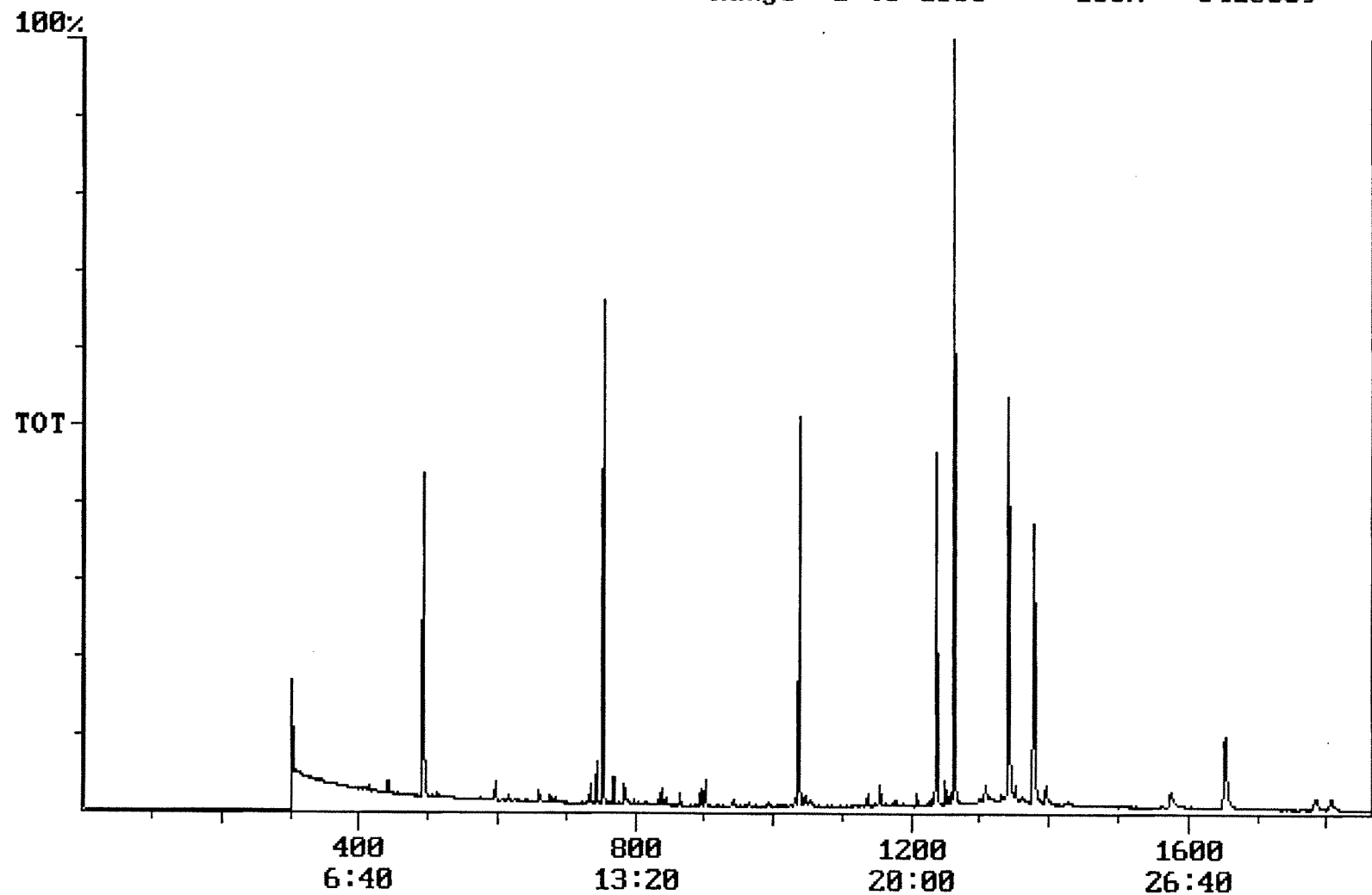


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/05/2001 Time: 10:19:56

Sample: AD22777
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		5.0		1.0
20	Surr_ Triphenyl phosphate		5.9		1.0
21	Surr_ Perylene-d12		4.3		1.0

Chromatogram Plot File: F:\22777FFF Date: Oct-02-1991 00:06:49
Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22777
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 6413859



Integration Report Quanfile: 22777FFF Quan Entries: 12
 Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22777
 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	1,482,002	1,005,927
2	phenanthrene d-10 (IS)	17:17	1037	Auto	2,211,664	1,090,243
3	chrysene d-12(IS)	22:58	1378	Auto	1,512,179	665,890
4	p-terphenyl d-14(IS)	21:02	1262	Auto	2,674,554	2,048,571
5	acenaphthene d-10(SURR)	12:33	753	Auto	1,482,002	1,005,927
6	phenanthrene d-10(SURR)	17:17	1037	Auto	2,211,664	1,090,243
7	chrysene d-12 (SURR)	22:58	1378	Auto	1,512,179	665,890
8	1,3-dimethyl-2-nitrobenzene	8:12	492	Auto	965,715	413,924
9	triphenyl phosphate (S)	22:20	1340	Auto	1,199,737	519,585
10	perylene d-12 (SURR)	27:33	1653	Auto	803,569	149,808
11	bis(2-ethylhexyl)phthalate	23:15	1395	Auto	161,430	41,448
12	2,6-dinitrotoluene	12:12	732	Auto	16,459	9,889

Quantitation Report

Quanfile: 22777FFF

Quan Entries: 12

Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22777

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	3.701	UG/L
6	phenanthrene d-10 (SURR)	A	1037	17:17	BV	3.602	UG/L
7	chrysene d-12 (SURR)	A	1378	22:58	BB	3.633	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	492	8:12	BB	4.997	UG/L
9	triphenyl phosphate (S)	A	1340	22:20	BB	5.892	UG/L
10	perylene d-12 (SURR)	A	1653	27:33	BB	4.278	UG/L
17	bis(2-ethylhexyl)phthalate	A	1395	23:15	BB	0.373	UG/L
20	2,6-dinitrotoluene	A	732	12:12	BB	0.160	UG/L
<i>reversed IS/SURR's</i>							

Comments for Sample AD22777 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 10-08-2001 Time: 13:05:38

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.