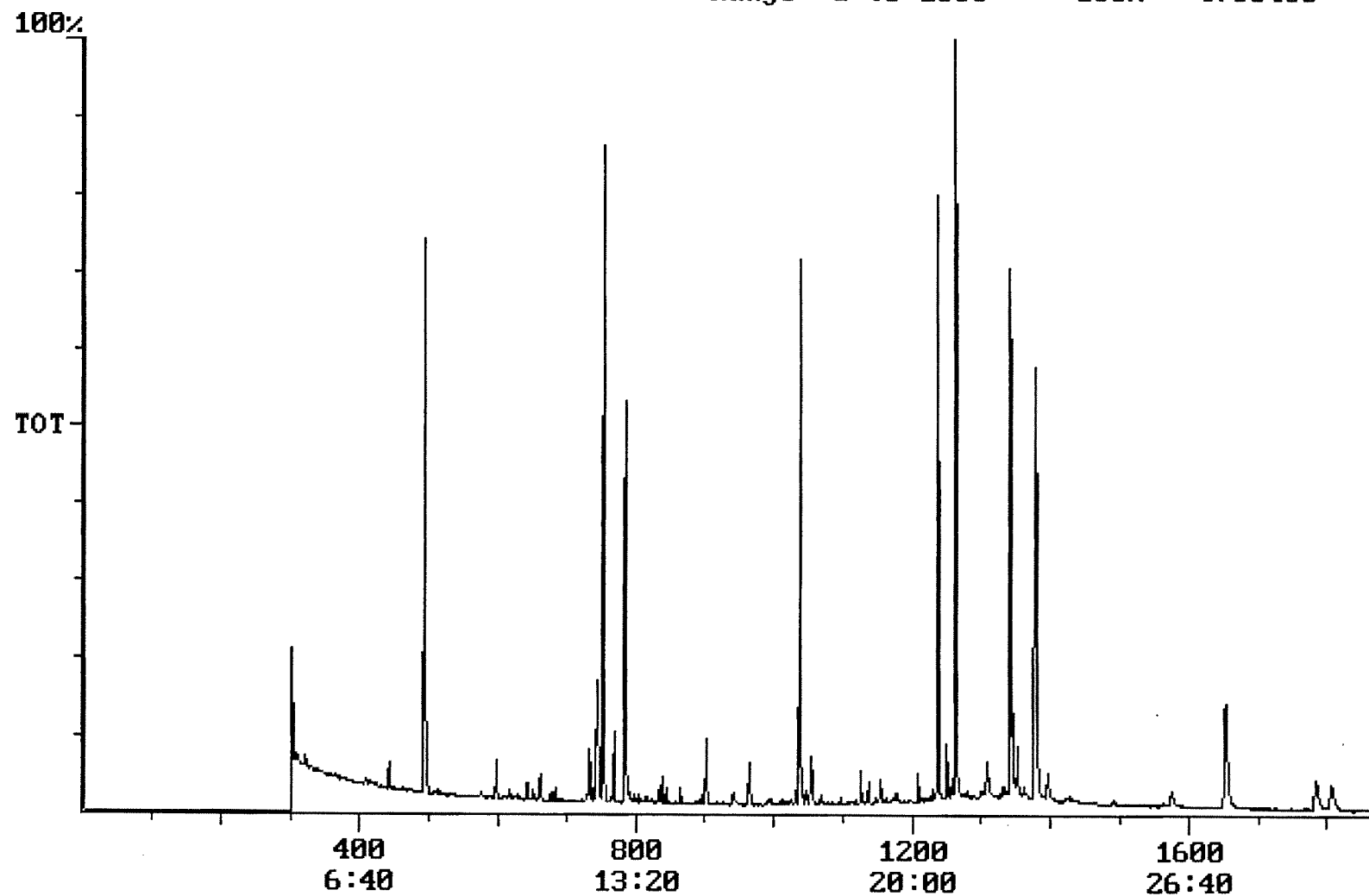


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/05/2001 Time: 10:21:18

Sample: AD22779
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		4.7		1.0
20	Surr_Triphenyl phosphate		5.9		1.0
21	Surr_Perylene-d12		4.2		1.0

Chromatogram Plot File: F:\22779HHH Date: Oct-02-1991 01:15:42
 Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22779
 Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
 Plotted: 1 to 1860 Range: 1 to 1860 100% = 4735403



Integration Report

Quanfile: 22779HHH

Quan Entries: 12

Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22779

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	1,585,362	992,401
2	phenanthrene d-10 (IS)	17:17	1037	Auto	2,332,311	1,139,962
3	chrysene d-12 (IS)	22:58	1378	Auto	1,575,155	706,760
4	p-terphenyl d-14 (IS)	21:02	1262	Auto	2,505,235	1,525,268
5	acenaphthene d-10 (SURR)	12:33	753	Auto	1,585,362	992,401
6	phenanthrene d-10 (SURR)	17:17	1037	Auto	2,332,311	1,139,962
7	chrysene d-12 (SURR)	22:58	1378	Auto	1,575,155	706,760
8	1,3-dimethyl-2-nitrobenzene	8:13	493	Auto	971,605	531,504
9	triphenyl phosphate (S)	22:20	1340	Auto	1,259,862	503,618
10	perylene d-12 (SURR)	27:33	1653	Auto	822,987	152,275
11	bis(2-ethylhexyl)phthalate	23:15	1395	Auto	172,678	43,456
12	2,4-dinitrotoluene	13:19	799	Auto	3,210	1,982

Quantitation Report Quanfile: 22779HHH Quan Entries: 12
 Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22779
 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	BU	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	4.227	UG/L
6	phenanthrene d-10(SURR)	A	1037	17:17	BU	4.055	UG/L
7	chrysene d-12 (SURR)	A	1378	22:58	BB	4.040	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	493	8:13	BB	4.700	UG/L
9	triphenyl phosphate (S)	A	1340	22:20	BB	5.940	UG/L
10	perylene d-12 (SURR)	A	1653	27:33	BB	4.206	UG/L
17	bis(2-ethylhexyl)phthalate	A	1395	23:15	BB	0.383	UG/L
21	2,4-dinitrotoluene	A	799	13:19	BB	0.023	UG/L
reversed IS/surr's							

Comments for Sample AD22779 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 10-08-2001 Time: 13:10:42

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.