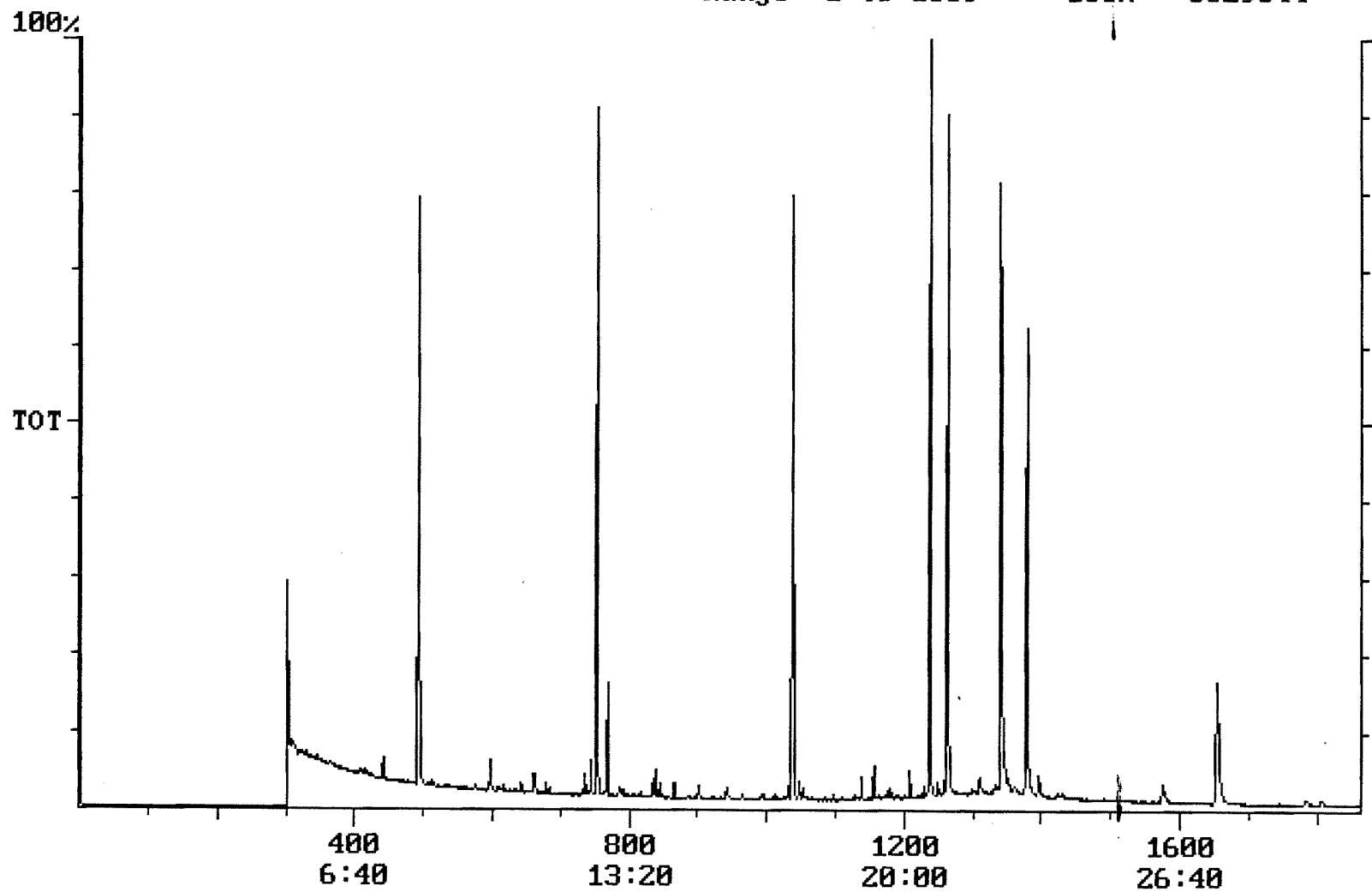


User: Bohlk, Ted
 Westchester Dept. of Labs & Research
 Date: 10/05/2001 Time: 10:18:25

Sample: AD22774
 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		4.9		1.0
20	Surr_Triphenyl phosphate		6.0		1.0
21	Surr_Perylene-d12		4.1		1.0

Chromatogram Plot File: F:\22774CCC Date: Oct-01-1991 22:23:32
Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22774
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1859 Range: 1 to 1859 100% = 3829344



Integration Report Quanfile: 22774CCC Quan Entries: 12
 Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22774
 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	1,279,983	801,250
2	phenanthrene d-10 (IS)	17:17	1037	Auto	2,107,123	1,032,846
3	chrysene d-12 (IS)	22:58	1378	Auto	1,506,845	677,909
4	p-terphenyl d-14 (IS)	21:02	1262	Auto	2,213,875	1,075,813
5	acenaphthene d-10 (SURR)	12:33	753	Auto	1,279,983	801,250
6	phenanthrene d-10 (SURR)	17:17	1037	Auto	2,107,123	1,032,846
7	chrysene d-12 (SURR)	22:58	1378	Auto	1,506,845	677,909
8	1,3-dimethyl-2-nitrobenzene	8:13	493	Auto	823,246	449,593
9	triphenyl phosphate (S)	22:20	1340	Auto	1,212,515	476,710
10	perylene d-12 (SURR)	27:33	1653	Auto	766,509	148,305
11	bis(2-ethylhexyl)phthalate	23:16	1396	Auto	125,304	27,096
12	2,6-dinitrotoluene	12:12	732	Auto	6,272	3,982

Quantitation Report

Quanfile: 22774CCC

Quan Entries: 12

Comment: BOHLK; METHOD 525; INST 204; 10/01/01; SAMPLE 22774

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	3.862	UG/L
6	phenanthrene d-10(SURR)	A	1037	17:17	BV	4.145	UG/L
7	chrysene d-12 (SURR)	A	1378	22:58	BB	4.373	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	493	8:13	BB	4.932	UG/L
9	triphenyl phosphate (S)	A	1340	22:20	BB	5.976	UG/L
10	perylene d-12 (SURR)	A	1653	27:33	BB	4.095	UG/L
17	bis(2-ethylhexyl)phthalate	A	1396	23:16	BB	0.291	UG/L
20	2,6-dinitrotoluene	A	732	12:12	BB	0.071	UG/L
<i>reversed IS/surr's</i>							

Comments for Sample AD22774 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 10-08-2001 Time: 12:56:57

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.