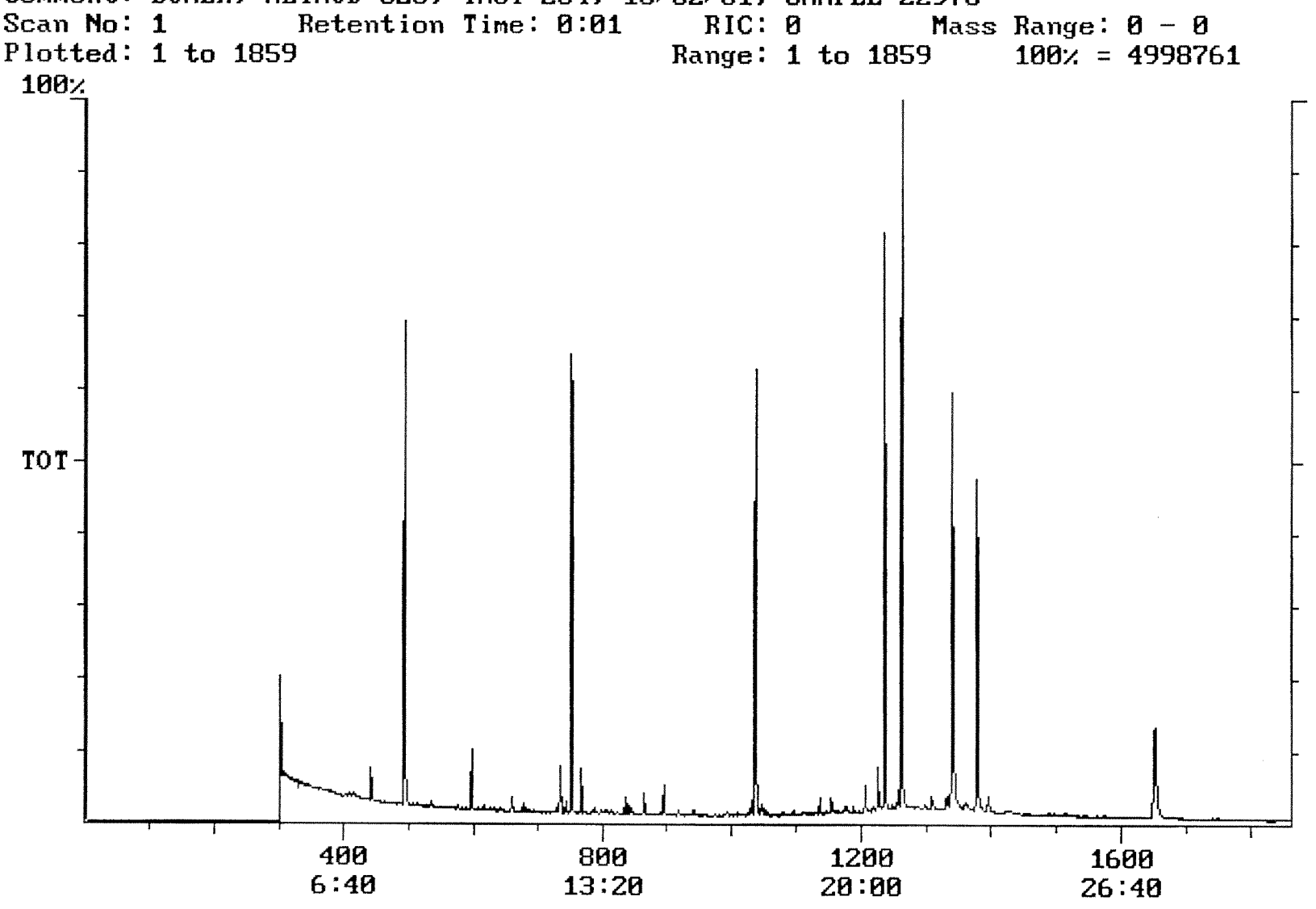
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/05/2001 Time: 16:30:06

Sample: AD22976
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 9/25/01 16:11 Ended: 10/2/01 16:11 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		6.1		1.0
21	Surr_Perylene-d12		4.3		1.0

Chromatogram Plot File: E:\22976J Date: Oct-03-1991 02:07:21
Comment: BOHLK; METHOD 525; INST 204; 10/02/01; SAMPLE 22976
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1859 Range: 1 to 1859 100% = 4998761



Integration Report Quanfile: 22976J Quan Entries: 12
Comment: BOHLK; METHOD 525; INST 204; 10/02/01; SAMPLE 22976
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:32	752	Auto	1,473,231	735,261
2	phenanthrene d-10 (IS)	17:16	1036	Auto	2,269,914	1,060,795
3	chrysene d-12(IS)	22:57	1377	Auto	1,541,103	665,463
4	p-terphenyl d-14(IS)	21:02	1262	Auto	2,314,528	1,478,779
5	acenaphthene d-10(SURR	12:32	752	Auto	1,473,231	735,261
6	phenanthrene d-10(SURR	17:16	1036	Auto	2,269,914	1,060,795
7	chrysene d-12 (SURR)	22:57	1377	Auto	1,541,103	665,463
8	1,3-dimethyl-2-nitrobe	8:12	492	Auto	903,959	517,500
9	triphenyl phosphate (S	22:20	1340	Auto	1,269,909	445,323
10	perylene d-12 (SURR)	27:31	1651	Auto	824,679	161,121
11	bis(2-ethylhexyl)phtha	23:15	1395	Auto	121,743	29,726
12	2,6-dinitrotoluene	12:11	731	Auto	17,224	10,314

Quantitation Report Quanfile: 22976J Quan Entries: 12
Comment: BOHLK; METHOD 525; INST 204; 10/02/01; SAMPLE 22976
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	752	12:32	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	1377	22:57	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	752	12:32	BB	4.251	UG/L
6	phenanthrene d-10(SURR	A	1036	17:16	BV	4.272	UG/L
7	chrysene d-12 (SURR)	A	1377	22:57	BB	4.278	UG/L
8	1,3-dimethyl-2-nitrobe	A	492	8:12	BB	4.706	UG/L
9	triphenyl phosphate (S	A	1340	22:20	BV	6.120	UG/L
10	perylene d-12 (SURR)	A	1651	27:31	BB	4.308	UG/L
17	bis(2-ethylhexyl)phtha	A	1395	23:15	BV	0.277	UG/L
20	2,6-dinitrotoluene	A	731	12:11	BB	0.169	UG/L

reviewed IS/ Surr's

Comments for Sample AD22976 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 10-08-2001 Time: 09:30:44

Recoveries for three of the eighteen compounds in the MDL Standard were above their acceptance ranges. The recovery for Terbacil in the LCS Standard was above its acceptance range. The recoveries for six of the eighteen compounds in the Spiked Sample were outside their acceptance ranges.