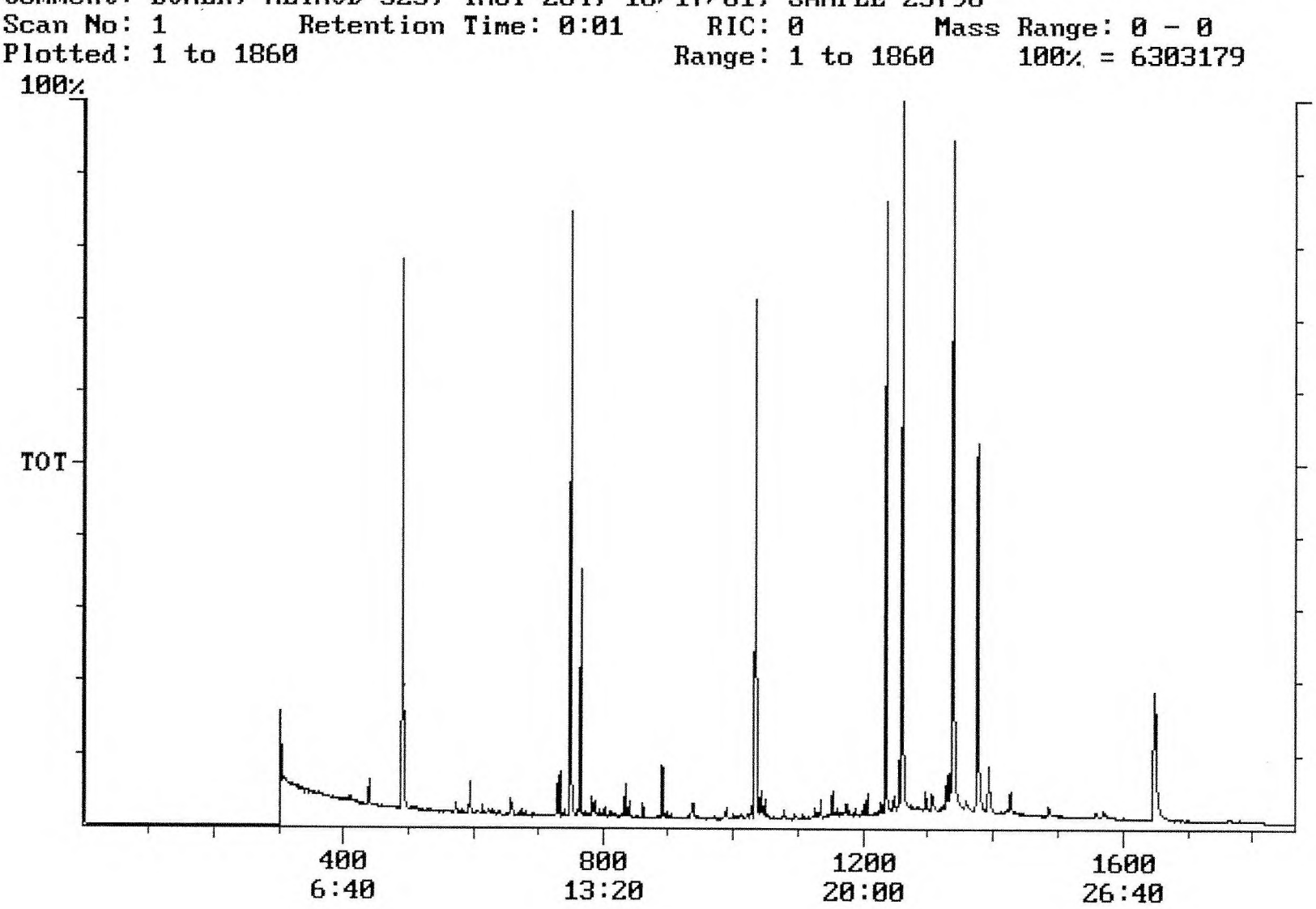
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 10/27/2001 Time: 15:37:37

Sample: AD23799
Analysis: S525 Organic Chemicals - 525.2
Units: ug
Started: 10/9/01 15:31 Ended: 10/17/01 15:31 Analyst: TB
Result source: manual entry
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	Hexachlorocyclopentadiene		< MDL		0.10
2	Hexachlorobenzene		< MDL		0.10
3	Simazine		< MDL		0.40
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-nitrobenze		5.1		1.0
20	Surr_Triphenyl phosphate		5.5		1.0
21	Surr_Perylene-d12		4.9		1.0

Chromatogram Plot File: F:\23799II Date: Oct-18-1991 02:37:34
Comment: BOHLK; METHOD 525; INST 204; 10/17/01; SAMPLE 23798
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1860 Range: 1 to 1860 100% = 6303179



Integration Report Quanfile: 23799II Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 10/17/01; SAMPLE 23798
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:30	750	Auto	2,167,012	1,237,352
2	phenanthrene d-10 (IS)	17:14	1034	Auto	3,327,806	1,535,013
3	chrysene d-12(IS)	22:55	1375	Auto	2,451,386	881,237
4	p-terphenyl d-14(IS)	21:00	1260	Auto	4,033,413	1,973,837
5	acenaphthene d-10(SURR	12:30	750	Auto	2,167,012	1,237,352
6	phenanthrene d-10(SURR	17:14	1034	Auto	3,327,806	1,535,013
7	chrysene d-12 (SURR)	22:55	1375	Auto	2,451,386	881,237
8	1,3-dimethyl-2-nitrobe	8:10	490	Auto	1,335,381	758,417
9	triphenyl phosphate (S	22:18	1338	Auto	1,938,187	877,307
10	perylene d-12 (SURR)	27:28	1648	Auto	1,496,641	296,991
11	bis(2-ethylhexyl)adipa	22:09	1329	Auto	43,708	8,754
12	bis(2-ethylhexyl)phtha	23:13	1393	Auto	382,739	109,138
13	2,6-dinitrotoluene	12:09	729	Auto	65,685	35,691
14	2,4-dinitrotoluene	13:17	797	Auto	2,851	1,471
15	4,4'-dde	20:56	1256	Auto	1,345	1,345

Quantitation Report Quanfile: 23799II Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 10/17/01; SAMPLE 23798
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	750	12:30	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1034	17:14	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1375	22:55	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1260	21:00	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	750	12:30	BB	3.857	UG/L
6	phenanthrene d-10(SURR	A	1034	17:14	BV	4.020	UG/L
7	chrysene d-12 (SURR)	A	1375	22:55	BB	4.280	UG/L
8	1,3-dimethyl-2-nitrobe	A	490	8:10	BV	5.143	UG/L
9	triphenyl phosphate (S	A	1338	22:18	BB	5.533	UG/L
10	perylene d-12 (SURR)	A	1648	27:28	BB	4.919	UG/L
16	bis(2-ethylhexyl)adipa	A	1329	22:09	MM	0.108	UG/L
17	bis(2-ethylhexyl)phtha	A	1393	23:13	VB	0.506	UG/L
20	2,6-dinitrotoluene	A	729	12:09	BB	0.430	UG/L
21	2,4-dinitrotoluene	A	797	13:17	BB	0.016	UG/L
28	4,4'-dde	A	1256	20:56	BB	0.005	UG/L

reviewed IS/Surr's

Comments for Sample AD23799 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 11-04-2001 Time: 10:39:12

Recoveries for three of the eighteen compounds in the MDL Standard and the LCS Standard were outside of their acceptance ranges. Recoveries for two of the eighteen compounds in the Spiked Sample and the Spiked Duplicate Sample were outside of their acceptance ranges.