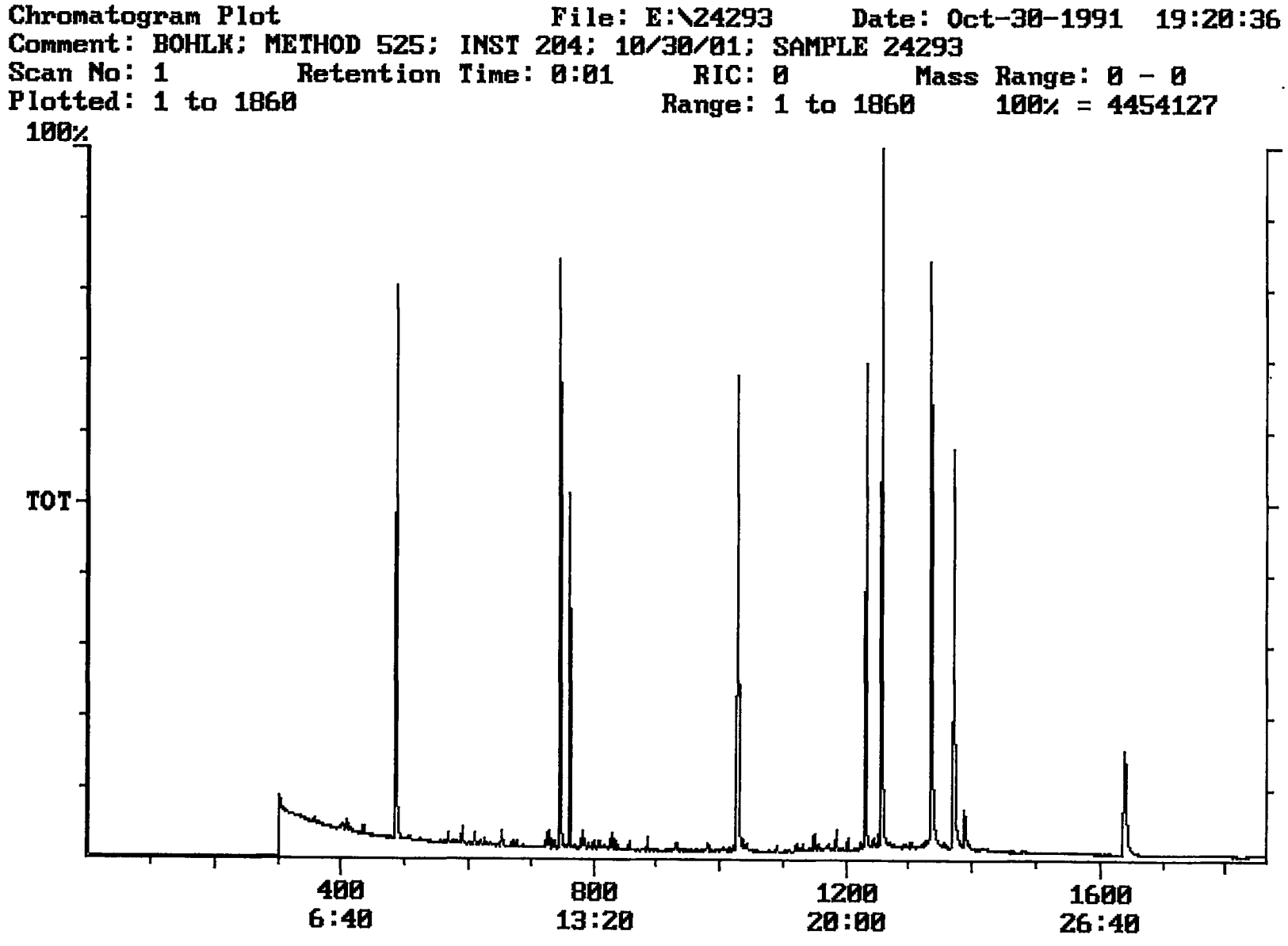
10/30
19:20

User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 11/13/2001 Time: 14:27:04

Sample: AD24293
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 10/16/01 14:24 Ended: 10/30/01 14:24 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		0.40
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-nitrobenzen		4.9		1.0
20	Surr_Triphenyl phosphate		5.6		1.0
21	Surr_Perylene-d12		4.3		1.0



Integration Report Quanfile: 24293 Quan Entries: 16
Comment: BOHLK; METHOD 525; INST 204; 10/30/01; SAMPLE 24293
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:26	746	Auto	1,573,002	894,392
2	phenanthrene d-10 (IS)	17:08	1028	Auto	2,448,499	1,024,631
3	chrysene d-12(IS)	22:51	1371	Auto	1,753,567	742,324
4	p-terphenyl d-14(IS)	20:57	1257	Auto	2,655,854	1,426,666
5	acenaphthene d-10(SURR	12:26	746	Auto	1,573,002	894,392
6	phenanthrene d-10(SURR	17:08	1028	Auto	2,448,499	1,024,631
7	chrysene d-12 (SURR)	22:51	1371	Auto	1,753,567	742,324
8	1,3-dimethyl-2-nitrobe	8:06	486	Auto	936,372	529,707
9	triphenyl phosphate (S	22:14	1334	Auto	1,337,810	554,968
10	perylene d-12 (SURR)	27:18	1638	Auto	951,908	161,714
11	hexachlorocyclopentadi	10:17	617	Auto	1,501	1,501
12	hexachlorobenzene	15:35	935	Auto	8,547	2,957
13	bis(2-ethylhexyl)adipa	22:14	1334	Auto	26,763	4,687
14	bis(2-ethylhexyl)phtha	23:08	1388	Auto	287,181	70,093
15	2,4-dinitrotoluene	13:13	793	Auto	1,054	1,054
16	4,4'-dde	20:53	1253	Auto	7,977	5,626

reviewed IS/ surr's

Quantitation Report Quanfile: 24293 Quan Entries: 16
Comment: BOHLK; METHOD 525; INST 204; 10/30/01; SAMPLE 24293
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	746	12:26	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1028	17:08	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1371	22:51	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1257	20:57	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	746	12:26	BB	4.687	UG/L
6	phenanthrene d-10(SURR	A	1028	17:08	BV	4.974	UG/L
7	chrysene d-12 (SURR)	A	1371	22:51	BB	4.883	UG/L
8	1,3-dimethyl-2-nitrobe	A	486	8:06	BV	4.869	UG/L
9	triphenyl phosphate (S	A	1334	22:14	BB	5.598	UG/L
10	perylene d-12 (SURR)	A	1638	27:18	BB	4.317	UG/L
11	hexachlorocyclopentadi	A	617	10:17	BB	0.023	UG/L
12	hexachlorobenzene	A	935	15:35	BB	0.041	UG/L
16	bis(2-ethylhexyl)adipa	A	1334	22:14	MM	0.084	UG/L
17	bis(2-ethylhexyl)phtha	A	1388	23:08	VB	0.524	UG/L
21	2,4-dinitrotoluene	A	793	13:13	BB	0.010	UG/L
28	4,4'-dde	A	1253	20:53	BB	0.037	UG/L

Comments for Sample AD24293 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 11-21-2001 Time: 16:24:17

The recoveries for 4 of the 18 compounds in the MDL Standard and 2 of the 18 compounds in the LCS Standard were outside of their acceptance ranges.