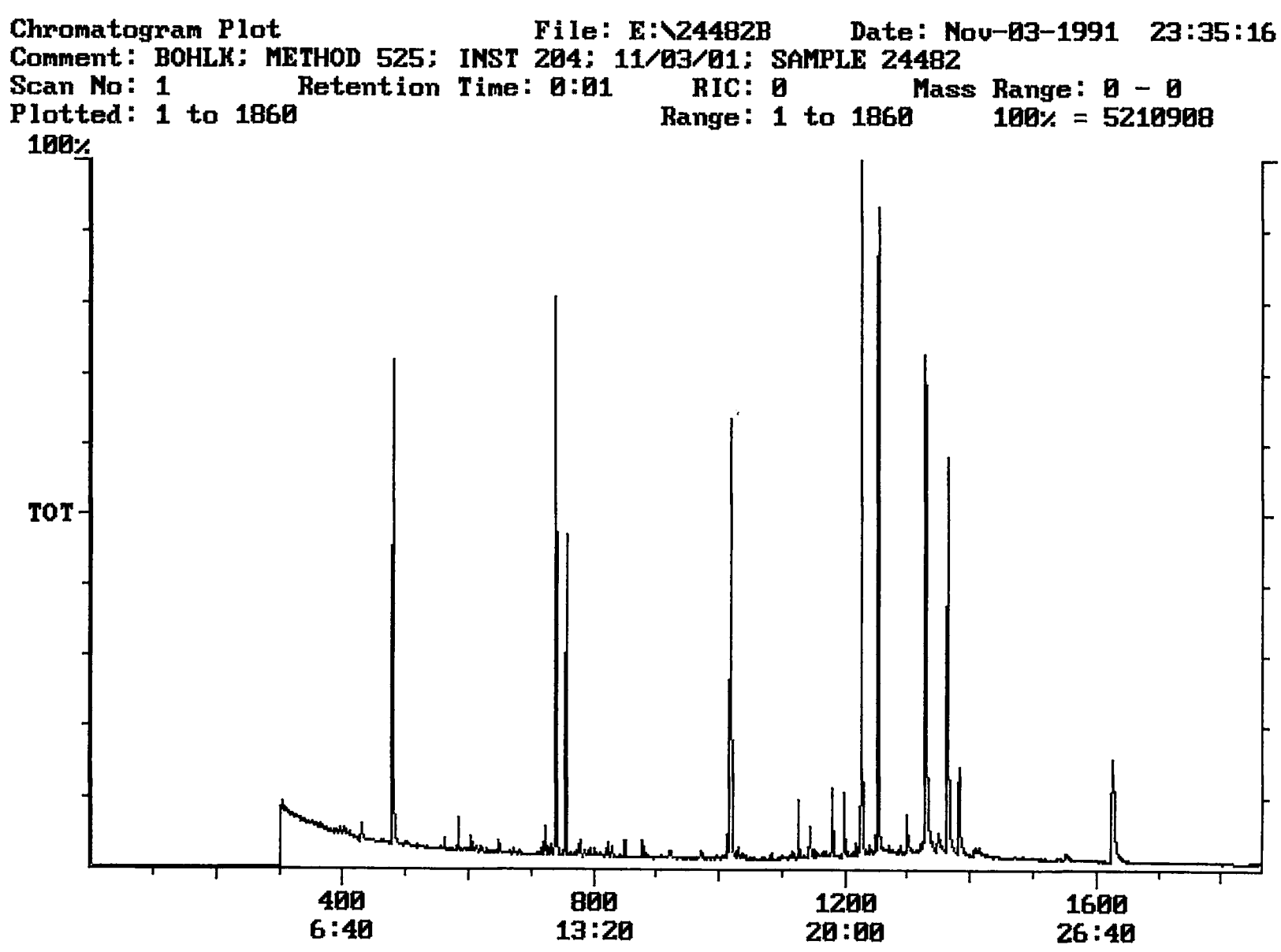
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 11/17/2001 Time: 14:47:02

Sample: AD24482
Analysis: S525 Organic Chemicals - 525.2
Units: ug
Started: 11/17/01 14:45 Ended: 11/17/01 14:45 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		1.1		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		5.2		1.0
20	Surr_Triphenyl phosphate		6.2		1.0
21	Surr_Perylene-d12		4.0		1.0



Integration Report Quanfile: 24482B Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/03/01; SAMPLE 24482
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:20	740	Auto	1,603,947	943,181
2	phenanthrene d-10 (IS)	16:59	1019	Auto	2,565,830	1,082,593
3	chrysene d-12(IS)	22:44	1364	Auto	1,939,934	860,507
4	p-terphenyl d-14(IS)	20:53	1253	Auto	2,909,275	1,282,000
5	acenaphthene d-10(SURR	12:20	740	Auto	1,603,947	943,181
6	phenanthrene d-10(SURR	16:59	1019	Auto	2,565,830	1,082,593
7	chrysene d-12 (SURR)	22:44	1364	Auto	1,939,934	860,507
8	1,3-dimethyl-2-nitrobe	8:00	480	Auto	1,029,093	570,796
9	triphenyl phosphate (S	22:08	1328	Auto	1,572,207	540,688
10	perylene d-12 (SURR)	27:06	1626	Auto	1,041,298	193,020
11	hexachlorobenzene	15:24	924	Auto	5,278	2,319
12	bis(2-ethylhexyl)adipa	22:09	1329	Auto	18,322	2,126
13	bis(2-ethylhexyl)phtha	23:02	1382	Auto	682,624	182,240
14	4,4'-dde	20:48	1248	Auto	4,553	3,477

reviewed IS/Surr's

Quantitation Report Quanfile: 24482B Quan Entries: 14
Comment: BOHLK; METHOD 525; INST 204; 11/03/01; SAMPLE 24482
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	740	12:20	BV	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1019	16:59	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1364	22:44	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1253	20:53	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	740	12:20	BV	3.698	UG/L
6	phenanthrene d-10(SURR	A	1019	16:59	BV	3.934	UG/L
7	chrysene d-12 (SURR)	A	1364	22:44	BB	4.131	UG/L
8	1,3-dimethyl-2-nitrobe	A	480	8:00	BV	5.169	UG/L
9	triphenyl phosphate (S	A	1328	22:08	BB	6.153	UG/L
10	perylene d-12 (SURR)	A	1626	27:06	BB	3.956	UG/L
12	hexachlorobenzene	A	924	15:24	BB	0.025	UG/L
16	bis(2-ethylhexyl)adipa	A	1329	22:09	MM	0.088	UG/L
17	bis(2-ethylhexyl)phtha	A	1382	23:02	VB	1.056	UG/L
28	4,4'-dde	A	1248	20:48	BB	0.017	UG/L

Comments for Sample AD24482 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 11-24-2001 Time: 12:23:05

Recovery for 1 of the 18 compounds in each of the MDL and LCS Standards was above its acceptance range.