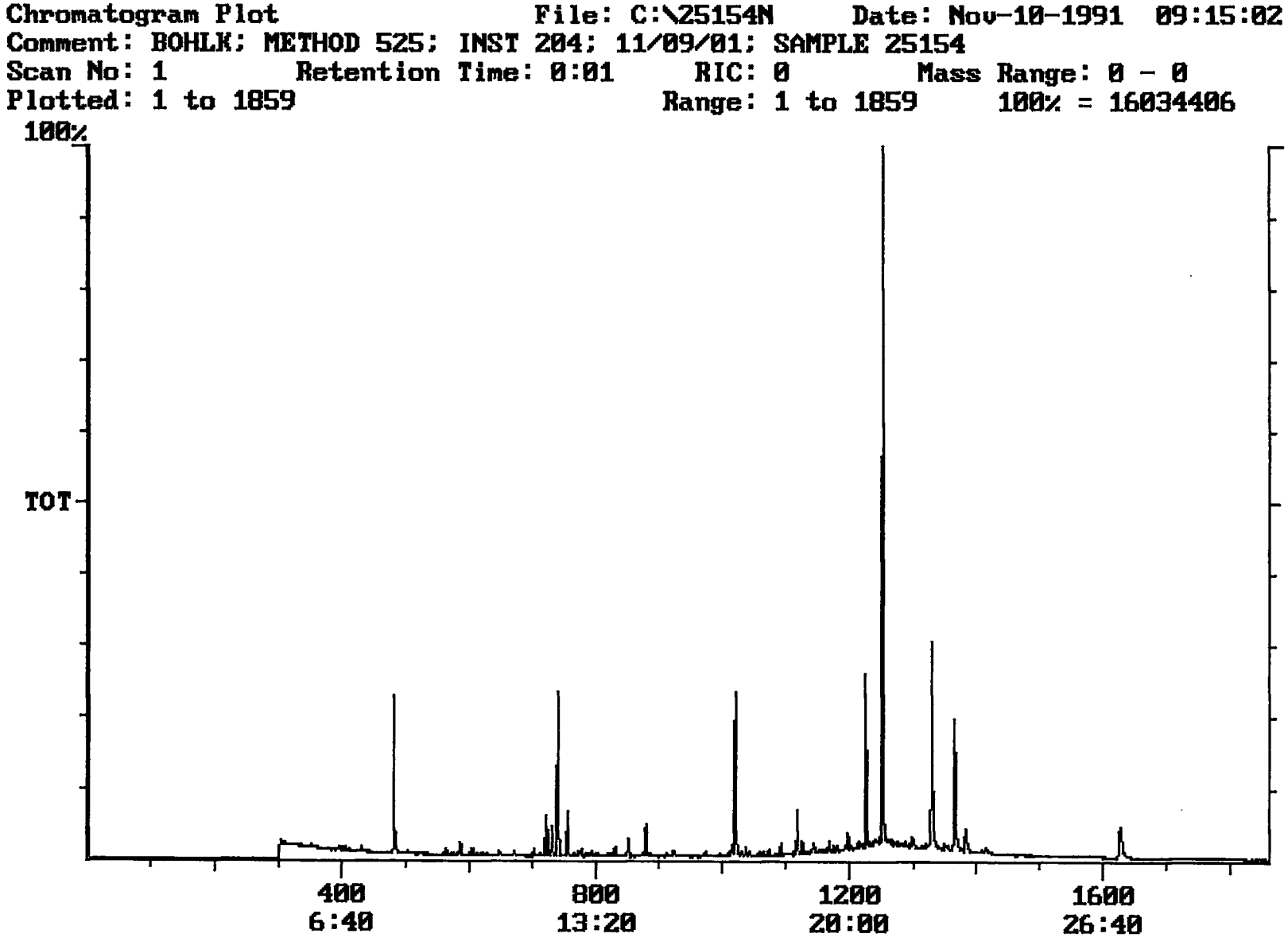
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 11/29/2001 Time: 10:23:07

Sample: AD25154
Analysis: S525 Organic Chemicals - 525.2
Units: ug
Started: 10/25/01 10:19 Ended: 11/9/01 10:19 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		0.20
5	Alachlor		< MDL		0.20
6	bis(2-Ethylhexyl)adipate		< MDL		0.60
7	bis(2-Ethylhexyl)phthalate		0.75		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		5.3		1.0
20	Surr_Triphenyl phosphate		5.7		1.0
21	Surr_Perylene-d12		3.4		1.0



Integration Report Quanfile: 25154N Quan Entries: 13
Comment: BOHLK; METHOD 525; INST 204; 11/09/01; SAMPLE 25154
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	1,759,357	849,318
2	phenanthrene d-10 (IS)	17:00	1020	Auto	2,828,441	1,209,132
3	chrysene d-12(IS)	22:45	1365	Auto	2,026,779	817,962
4	p-terphenyl d-14(IS)	20:53	1253	Auto	7,198,364	5,501,453
5	acenaphthene d-10(SURR	12:20	740	Auto	1,759,357	849,318
6	phenanthrene d-10(SURR	17:00	1020	Auto	2,828,441	1,209,132
7	chrysene d-12 (SURR)	22:45	1365	Auto	2,026,779	817,962
8	1,3-dimethyl-2-nitrobe	8:01	481	Auto	1,150,685	577,802
9	triphenyl phosphate (S	22:09	1329	Auto	1,518,123	705,689
10	perylene d-12 (SURR)	27:07	1627	Auto	938,077	171,680
11	bis(2-ethylhexyl)adipa	22:09	1329	Auto	35,769	5,201
12	bis(2-ethylhexyl)phtha	23:02	1382	Auto	508,136	142,969
13	2,6-dinitrotoluene	12:00	720	Auto	80,680	47,560

reviewed IS/surr's

Quantitation Report Quanfile: 25154N Quan Entries: 13
Comment: BOHLK; METHOD 525; INST 204; 11/09/01; SAMPLE 25154
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	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1020	17:00	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1365	22:45	BV	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1253	20:53	BV	5.000	UG/L
5	acenaphthene d-10(SURR	A	740	12:20	BB	1.640	UG/L
6	phenanthrene d-10(SURR	A	1020	17:00	BV	1.753	UG/L
7	chrysene d-12 (SURR)	A	1365	22:45	BV	1.745	UG/L
8	1,3-dimethyl-2-nitrobe	A	481	8:01	BV	5.269	UG/L
9	triphenyl phosphate (S	A	1329	22:09	BB	5.687	UG/L
10	perylene d-12 (SURR)	A	1627	27:07	BB	3.411	UG/L
16	bis(2-ethylhexyl)adipa	A	1329	22:09	MM	0.164	UG/L
17	bis(2-ethylhexyl)phtha	A	1382	23:02	BB	0.752	UG/L
20	2,6-dinitrotoluene	A	720	12:00	BB	0.703	UG/L

Comments for Sample AD25154 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-06-2001 Time: 11:16:46

Recoveries for 3 of the 18 compounds in the MDL and LCS Standards were outside of their acceptance ranges. Recoveries for 4 of the 18 compounds in the Spiked Sample were below their acceptance ranges.