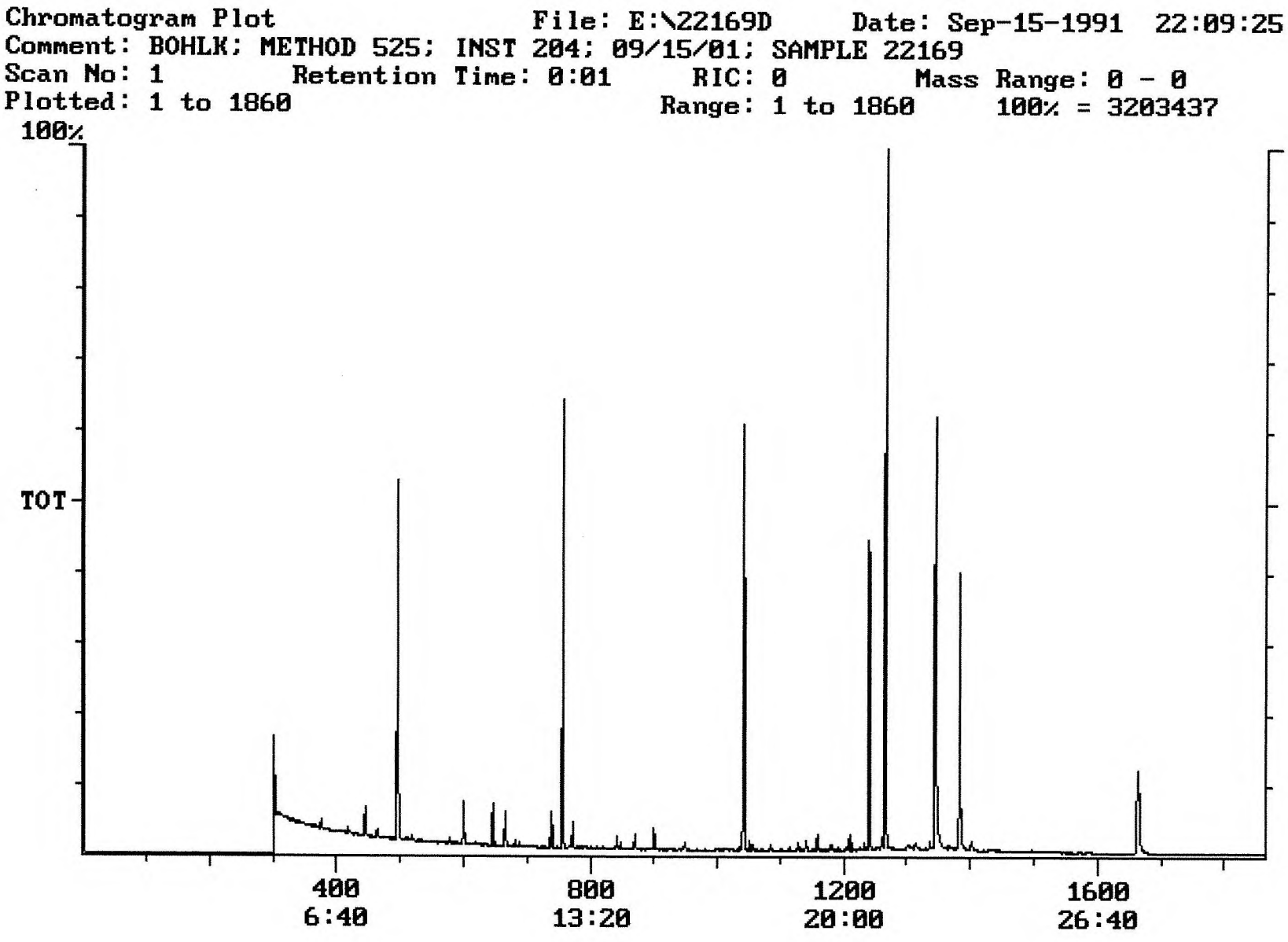
User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 09/17/2001 Time: 17:46:27

Sample: AD22169
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 9/12/01 17:41 Ended: 9/15/01 17:41 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		5.1		
20	Surr_Triphenyl phosphate		5.9		
21	Surr_Perylene-d12		4.7		



NYC 9/11 Public Portal Document

NYC-WTC_000063002

Integration Report Quanfile: 22169D Quan Entries: 13
Comment: BOHLK; METHOD 525; INST 204; 09/15/01; SAMPLE 22169
Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10(IS)	12:37	757	Auto	779,975	476,886
2	phenanthrene d-10 (IS)	17:22	1042	Auto	1,258,104	631,831
3	chrysene d-12(IS)	23:03	1383	Auto	885,454	368,647
4	p-terphenyl d-14(IS)	21:06	1266	Auto	1,452,651	1,094,626
5	acenaphthene d-10(SURR	12:37	757	Auto	779,975	476,886
6	phenanthrene d-10(SURR	17:22	1042	Auto	1,258,104	631,831
7	chrysene d-12 (SURR)	23:03	1383	Auto	885,454	368,647
8	1,3-dimethyl-2-nitrobe	8:16	496	Auto	487,694	254,450
9	triphenyl phosphate (S	22:25	1345	Auto	730,797	311,968
10	perylene d-12 (SURR)	27:43	1663	Auto	534,402	101,068
11	bis(2-ethylhexyl)phtha	23:21	1401	Auto	61,800	13,836
12	2,6-dinitrotoluene	12:15	735	Auto	14,647	7,805
13	2,4-dinitrotoluene	13:22	802	Auto	1,000	1,000

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NYC-WTC_000063003

Quantitation Report Quanfile: 22169D Quan Entries: 13
Comment: BOHLK; METHOD 525; INST 204; 09/15/01; SAMPLE 22169
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	757	12:37	BB	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1042	17:22	BB	5.000	UG/L
3	chrysene d-12(IS)	I	1383	23:03	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1266	21:06	BB	5.000	UG/L
5	acenaphthene d-10(SURR	A	757	12:37	BB	3.715	UG/L
6	phenanthrene d-10(SURR	A	1042	17:22	BB	4.035	UG/L
7	chrysene d-12 (SURR)	A	1383	23:03	BB	3.969	UG/L
8	1,3-dimethyl-2-nitrobe	A	496	8:16	BB	5.105	UG/L
9	triphenyl phosphate (S	A	1345	22:25	BB	5.898	UG/L
10	perylene d-12 (SURR)	A	1663	27:43	BB	4.702	UG/L
17	bis(2-ethylhexyl)phtha	A	1401	23:21	BB	0.309	UG/L
20	2,6-dinitrotoluene	A	735	12:15	BB	0.302	UG/L
21	2,4-dinitrotoluene	A	802	13:22	BB	0.016	UG/L

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Comments for Sample AD22169 - Analysis $525

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 09-18-2001 Time: 08:44:00

The recovery for Terbacil in the LCS Standard was above its acceptance range.