

User: Bohlk, Ted
 Westchester Dept. of Labs & Research
 Date: 09/25/2001 Time: 17:26:55

Sample: AD22492
 Analysis: \$525 Organic Chemicals - 525.2
 Units: ug
 Started: 9/19/01 17:14 Ended: 9/21/01 17:14 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		1.0
20	Surr_Triphenyl phosphate *		9.0		1.0
21	Surr_Perylene-d12		3.6		1.0

← flag T.V. = 5.0

Chromatogram Plot

File: E:\22492F

Date: Sep-21-1991 21:09:34

Comment: BOHLK; METHOD 525; INST 204; 09/21/01; SAMPLE 22492

Scan No: 1

Retention Time: 0:01

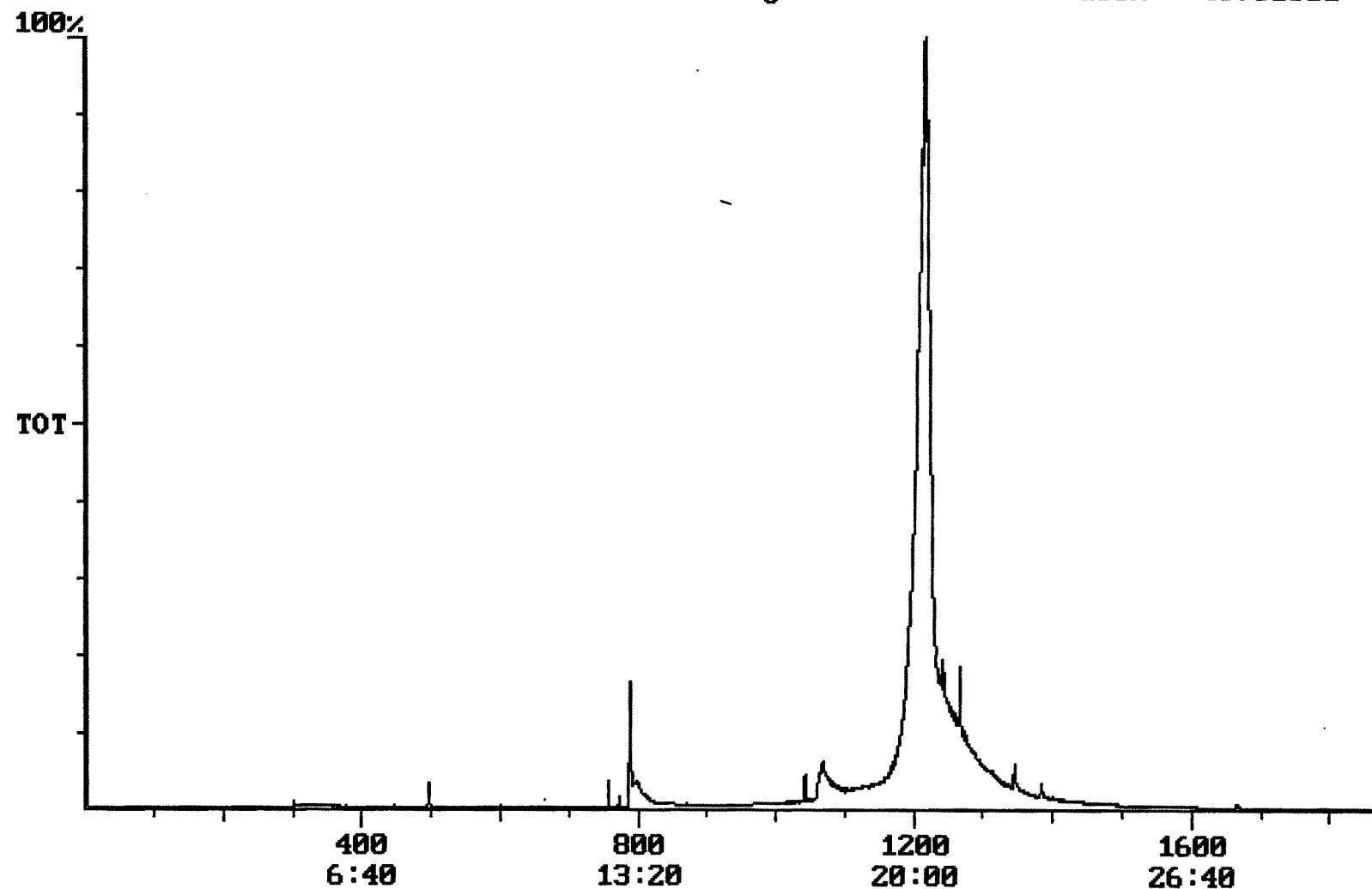
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1860

Range: 1 to 1860

100% = 45761311



Integration Report

Quanfile: 22492F

Quan Entries: 12

Comment: BOHLK; METHOD 525; INST 204; 09/21/01; SAMPLE 22492

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	623,652	381,735
2	phenanthrene d-10 (IS)	17:23	1043	Auto	1,222,087	516,114
3	chrysene d-12 (IS)	23:04	1384	Auto	581,419	233,071
4	p-terphenyl d-14 (IS)	21:07	1267	Auto	2,918,125	1,865,697
5	acenaphthene d-10 (SURR)	12:37	757	Auto	623,652	381,735
6	phenanthrene d-10 (SURR)	17:23	1043	Auto	1,222,087	516,114
7	chrysene d-12 (SURR)	23:04	1384	Auto	581,419	233,071
8	1,3-dimethyl-2-nitrobenzene	8:16	496	Auto	390,599	225,436
9	triphenyl phosphate (S)	22:25	1345	Auto	731,239	253,571
10	perylene d-12 (SURR)	27:45	1665	Auto	269,235	49,947
11	bis(2-ethylhexyl)phthalate	23:22	1402	Auto	58,566	14,865
12	2,6-dinitrotoluene	12:16	736	Auto	10,186	6,449

Quantitation Report Quanfile: 22492F Quan Entries: 12
 Comment: BOHLK; METHOD 525; INST 204; 09/21/01; SAMPLE 22492
 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	1043	17:23	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1384	23:04	BB	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1267	21:07	BB	5.000	UG/L
5	acenaphthene d-10(SURR)	A	757	12:37	BB	1.479	UG/L
6	phenanthrene d-10(SURR)	A	1043	17:23	BV	1.951	UG/L
7	chrysene d-12 (SURR)	A	1384	23:04	BB	1.298	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	496	8:16	BB	5.114	UG/L
9	triphenyl phosphate (S)	A	1345	22:25	BB	8.987	UG/L
10	perylene d-12 (SURR)	A	1665	27:45	BB	3.608	UG/L
17	bis(2-ethylhexyl)phthalate	A	1402	23:22	MM	0.442	UG/L
20	2,6-dinitrotoluene	A	736	12:16	BB	0.263	UG/L

Comments for Sample AD22492 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 09-26-2001 Time: 15:17:07

The recovery for one of the three Surrogate Standards in this sample was above its acceptance range. The recovery for Terbacil in the LCS Standard was above its acceptance range.