User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 12/21/2001 Time: 16:23:19

Sample: AD26711
Analysis: $525 Organic Chemicals - 525.2
Units: ug
Started: 11/8/01 16:21 Ended: 11/28/01 16:21 Analyst: TB
Result source: manual entry
Page: 1

| | Component Name | Viol | Result | 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 | | 0.82 | 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 | | 5.3 | 1.0 |
| 20 | Surr_Triphenyl phosphate | USPEC | 6.8 | 1.0 |
| 21 | Surr_Perylene-d12 | | 5.7 | 1.0 |

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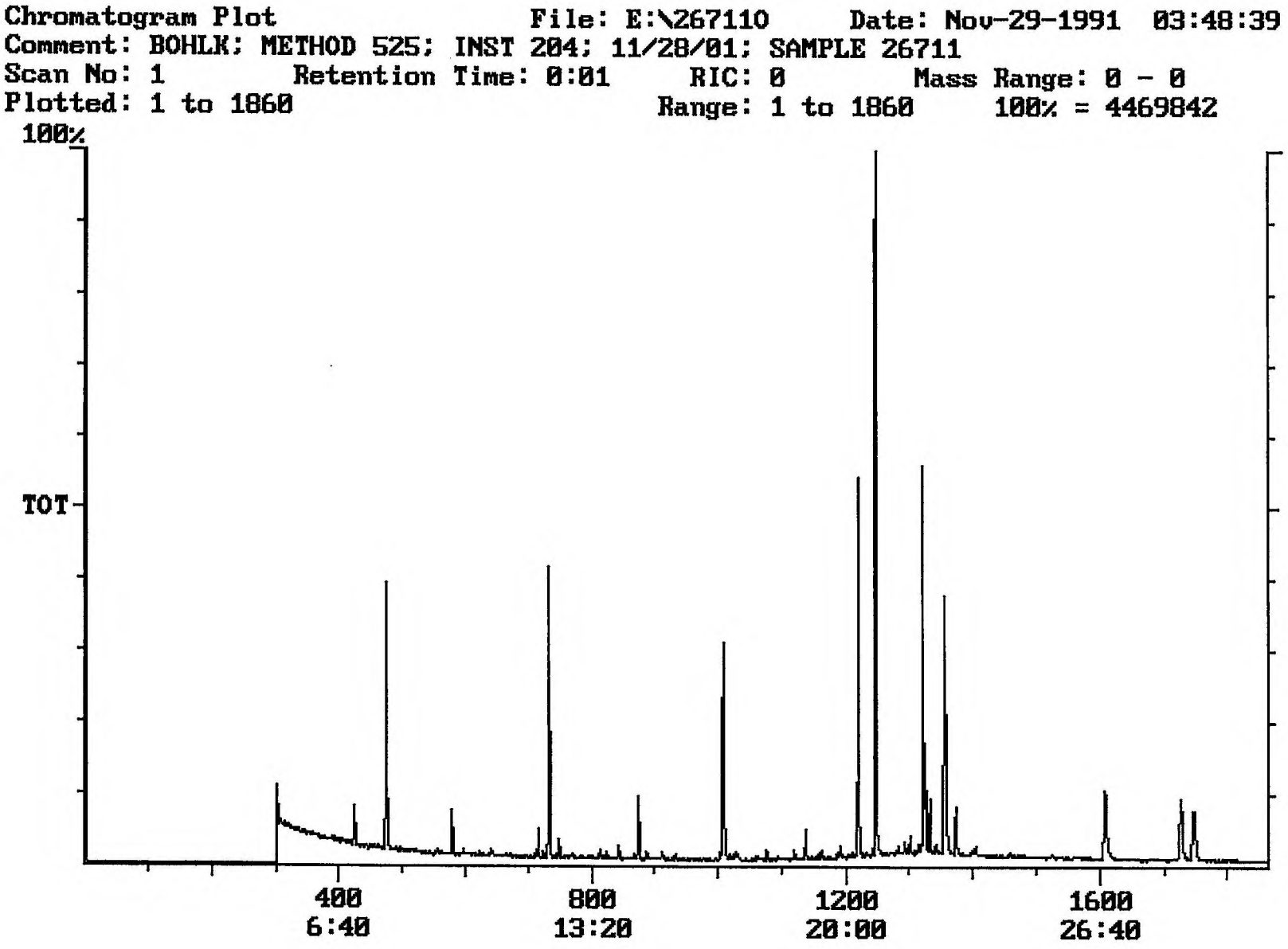

NYC-WTC_000030175

NYC 9/11 Public Portal Document

Integration Report  Quanfile: 267110  Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26711
Sorted via: Entry Number ↑

| No | Name of Compound | R Time | Scan# | Mode | Peak Area | Pk Height |
|---|---|---|---|---|---|---|
| 1 | acenaphthene d-10(IS) | 12:12 | 732 | Auto | 806,082 | 405,458 |
| 2 | phenanthrene d-10 (IS) | 16:48 | 1008 | Auto | 1,200,948 | 457,816 |
| 3 | chrysene d-12(IS) | 22:36 | 1356 | Auto | 838,694 | 278,430 |
| 4 | p-terphenyl d-14(IS) | 20:47 | 1247 | Auto | 2,728,263 | 1,273,697 |
| 5 | acenaphthene d-10(SURR | 12:12 | 732 | Auto | 806,082 | 405,458 |
| 6 | phenanthrene d-10(SURR | 16:48 | 1008 | Auto | 1,200,948 | 457,816 |
| 7 | chrysene d-12 (SURR) | 22:36 | 1356 | Auto | 838,694 | 278,430 |
| 8 | 1,3-dimethyl-2-nitrobe | 7:54 | 474 | Auto | 511,606 | 293,095 |
| 9 | triphenyl phosphate (S | 22:01 | 1321 | Auto | 632,220 | 300,939 |
| 10 | perylene d-12 (SURR) | 26:48 | 1608 | Auto | 554,712 | 104,233 |
| 11 | simazine | 16:01 | 961 | Auto | 0 | 0 |
| 12 | bis(2-ethylhexyl)adipa | 22:01 | 1321 | Auto | 28,054 | 7,048 |
| 13 | bis(2-ethylhexyl)phtha | 22:53 | 1373 | Auto | 249,369 | 90,931 |
| 14 | 2,6-dinitrotoluene | 11:53 | 713 | Auto | 12,919 | 6,847 |
| 15 | butachlor | 20:20 | 1220 | Auto | 1,027 | 1,027 |

renewed IS/SURR's

NYC-WTC_000030176

NYC 9/11 Public Portal Document

Quantitation Report  Quanfile: 267110  Quan Entries: 15
Comment: BOHLK; METHOD 525; INST 204; 11/28/01; SAMPLE 26711
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 | 732 | 12:12 | BB | 5.000 | UG/L |
| 2 | phenanthrene d-10 (IS) | I | 1008 | 16:48 | BV | 5.000 | UG/L |
| 3 | chrysene d-12(IS) | I | 1356 | 22:36 | BV | 5.000 | UG/L |
| 4 | p-terphenyl d-14(IS) | I | 1247 | 20:47 | BV | 5.000 | UG/L |
| 5 | acenaphthene d-10(SURR | A | 732 | 12:12 | BB | 1.949 | UG/L |
| 6 | phenanthrene d-10(SURR | A | 1008 | 16:48 | BV | 2.021 | UG/L |
| 7 | chrysene d-12 (SURR) | A | 1356 | 22:36 | BV | 1.837 | UG/L |
| 8 | 1,3-dimethyl-2-nitrobe | A | 474 | 7:54 | BV | 5.290 | UG/L |
| 9 | triphenyl phosphate (S | A | 1321 | 22:01 | BB | 6.800 | UG/L |
| 10 | perylene d-12 (SURR) | A | 1608 | 26:48 | BB | 5.729 | UG/L |
| 13 | simazine | A | 961 | 16:01 | BB | 0.001 | UG/L |
| 16 | bis(2-ethylhexyl)adipa | A | 1321 | 22:01 | MM | 0.149 | UG/L |
| 17 | bis(2-ethylhexyl)phtha | A | 1373 | 22:53 | BV | 0.817 | UG/L |
| 20 | 2,6-dinitrotoluene | A | 713 | 11:53 | BB | 0.275 | UG/L |
| 27 | butachlor | A | 1220 | 20:20 | BB | 0.012 | UG/L |

**Comments for Sample AD26711 - Analysis $525**

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 12-22-2001 Time: 13:03:41

The breakdown for Endrin in the Breakdown Standard was 40%. Recoveries for 4 of the 18 compounds in the MDL Standard were outside of their acceptance ranges. Recoveries for 4 of the 18 compounds in the LCS Standard were below their acceptance ranges. The recovery for 1 of the 3 Surrogate Standards in this sample was above its acceptance range.