

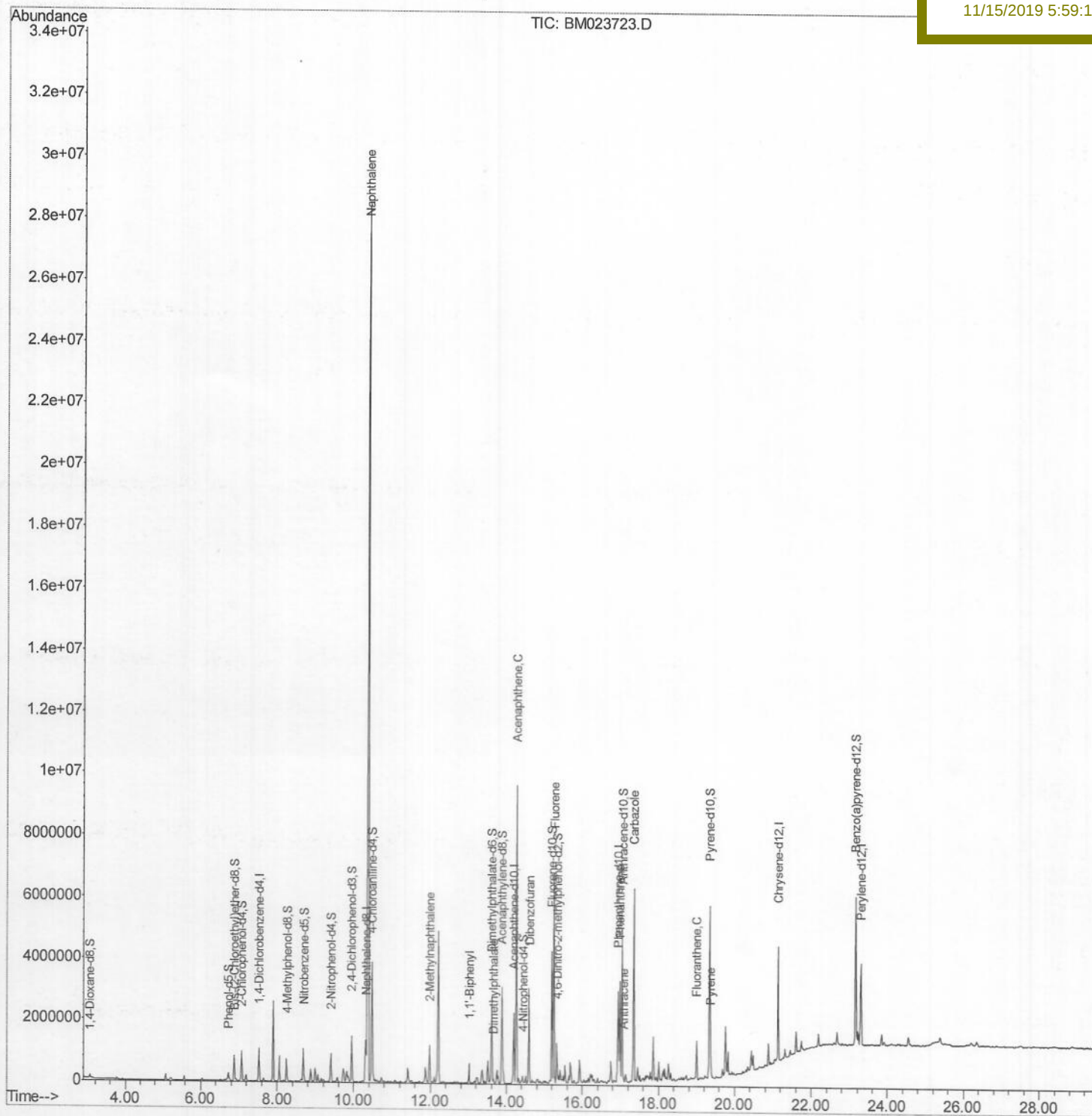
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111419\  
Data File : BM023723.D  
Acq On : 14 Nov 2019 19:05  
Operator : JU  
Sample : K5700-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 15 07:00:23 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM11119MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Nov 15 06:30:14 2019  
Response via : Initial Calibration

Instrument :  
BNA\_M  
ClientSampleId :  
BEGX1

Manual Integrations  
APPROVED

mohammad  
11/15/2019 5:59:19 PM



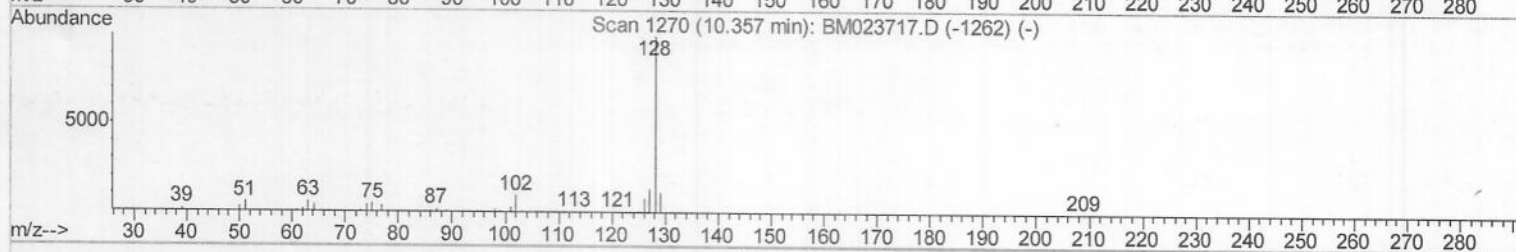
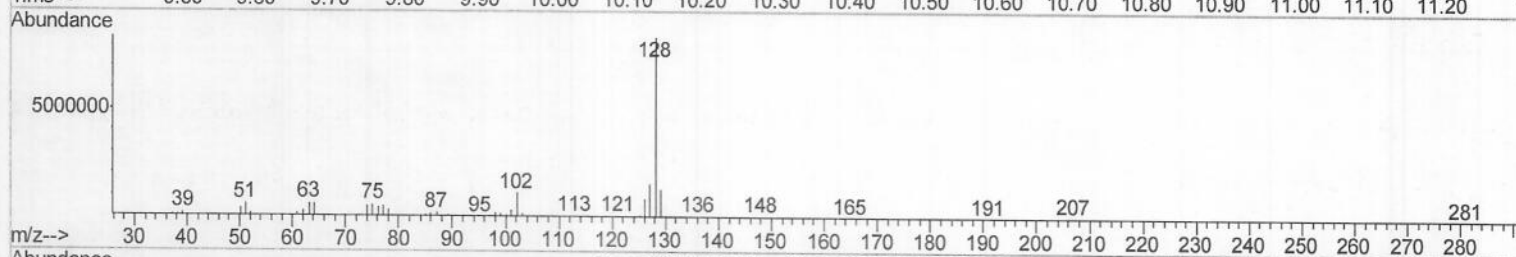
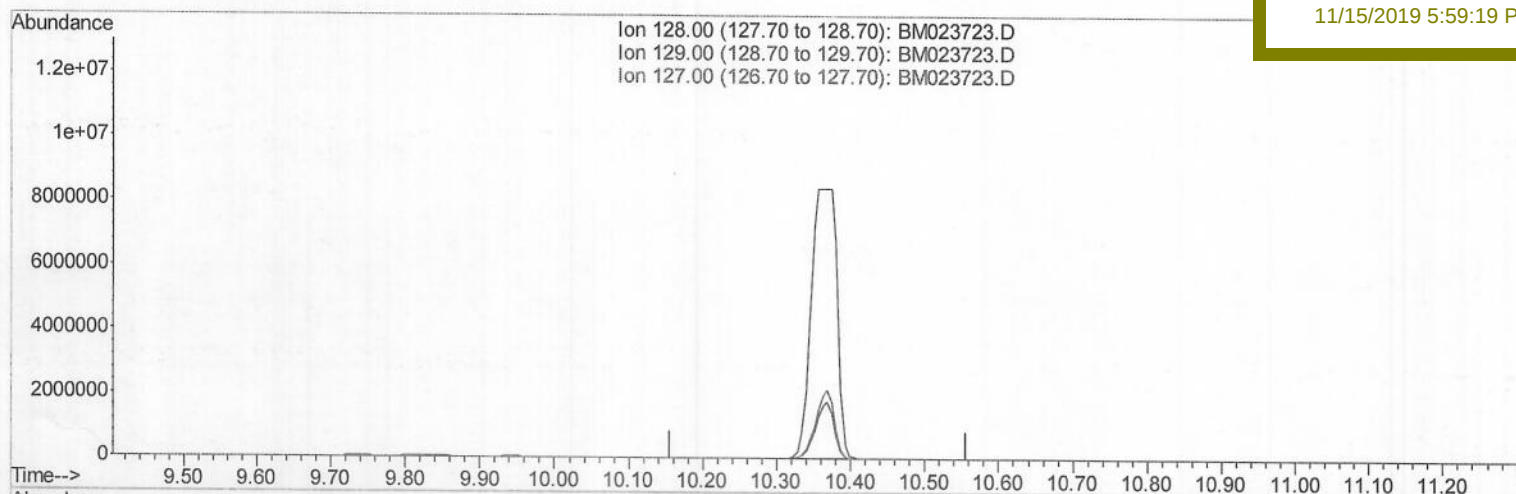
Data File : BM023723.D  
Acq On : 14 Nov 2019 19:05  
Operator : JU  
Sample : K5700-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 15 06:33:26 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM111119MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Nov 15 06:30:14 2019  
Response via : Initial Calibration

Instrument :  
BNA\_M  
ClientSampleId :  
BEGX1

Manual Integrations  
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TIC: BM023723.D

(28) Naphthalene

10.357min (-10.357) 0.00ng/ul

response 0

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 128.00 | 100   | 0.00  |
| 129.00 | 10.90 | 0.00# |
| 127.00 | 13.30 | 0.00# |
| 0.00   | 0.00  | 0.00  |

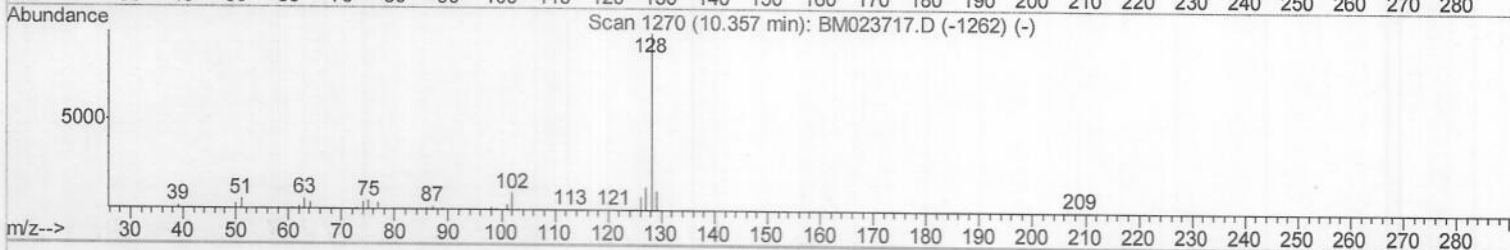
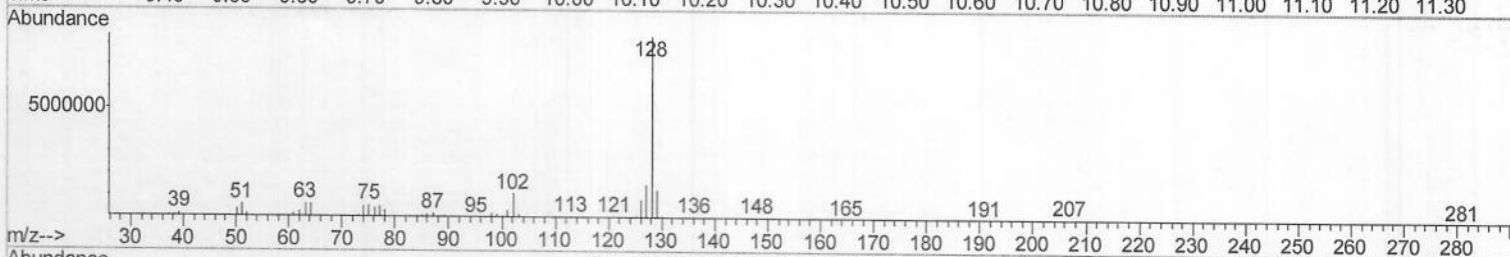
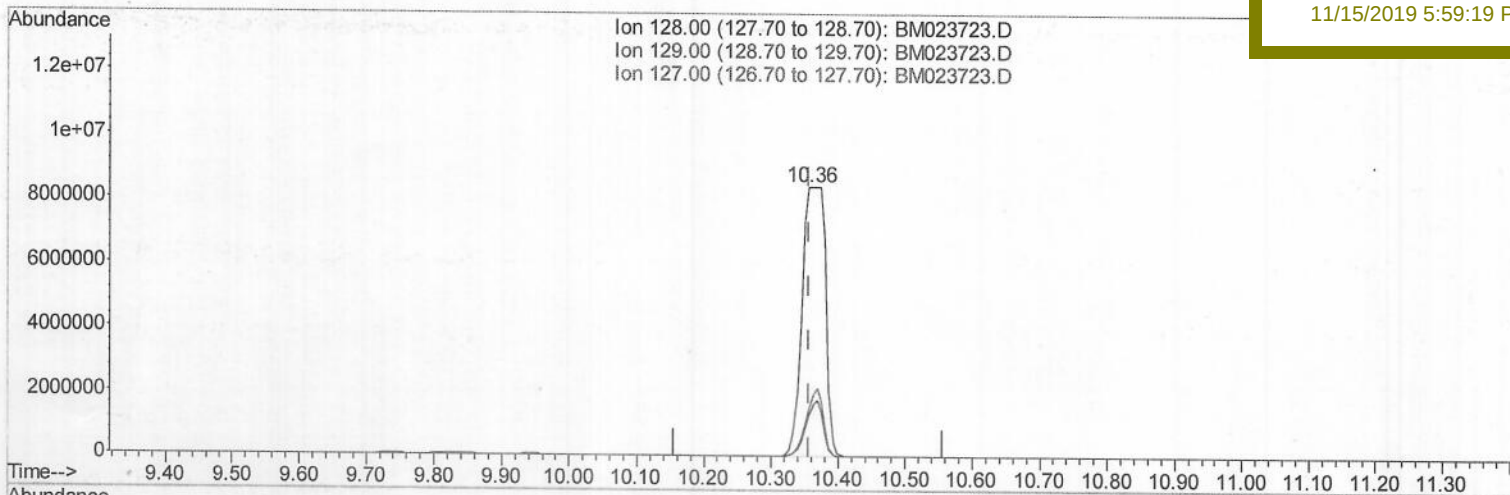
Data File : BM023723.D  
Acq On : 14 Nov 2019 19:05  
Operator : JU  
Sample : K5700-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 15 06:33:26 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM111119MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Nov 15 06:30:14 2019  
Response via : Initial Calibration

Instrument :  
BNA\_M  
ClientSampleId :  
BEGX1

Manual Integrations  
APPROVED

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11/15/2019 5:59:19 PM



TIC: BM023723.D

(28) Naphthalene

10.357min (+0.000) 318.39ng/ul m

JU 11/16/19

response 20295318

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 128.00 | 100   | 100    |
| 129.00 | 10.90 | 15.25# |
| 127.00 | 13.30 | 17.92# |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111419\  
 Data File : BM023723.D  
 Acq On : 14 Nov 2019 19:05  
 Operator : JU  
 Sample : K5700-01  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BEGX1

Quant Time: Nov 15 07:00:23 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM111119MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 15 06:30:14 2019  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

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 11/15/2019 5:59:19 PM

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.53  | 152  | 294683   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.31 | 136  | 1212777  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.19 | 164  | 769436   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.94 | 188  | 1659322  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.14 | 240  | 1632330  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.30 | 264  | 1923345  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.06  | 96  | 46344   | 6.48  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.72  | 99  | 153694  | 5.72  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl) ether-d | 6.87  | 67  | 378153  | 24.50 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.07  | 132 | 440908  | 22.66 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.25  | 113 | 283941  | 13.25 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.68  | 128 | 270833  | 31.20 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.40  | 143 | 310034  | 33.98 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.94  | 165 | 551383  | 26.50 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.46 | 131 | 58819   | 2.57  | ng/ul | 0.01 |
| 44) Dimethylphthalate-d6       | 13.61 | 166 | 1913473 | 31.89 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.87 | 160 | 2115372 | 30.89 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.42 | 143 | 61079   | 5.96  | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.19 | 176 | 1631395 | 30.79 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.32 | 200 | 295010  | 30.74 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.04 | 188 | 2659759 | 33.79 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.34 | 212 | 2902068 | 34.10 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.16 | 264 | 3419479 | 33.79 | ng/ul | 0.00 |

## Target Compounds

|                         |       |     |           |         |       |        |
|-------------------------|-------|-----|-----------|---------|-------|--------|
| 28) Naphthalene         | 10.36 | 128 | 20295318m | 318.389 | ng/ul | Ovalue |
| 34) 2-Methylnaphthalene | 11.98 | 142 | 528573    | 10.840  | ng/ul | 99     |
| 41) 1,1'-Biphenyl       | 13.02 | 154 | 330743    | 5.572   | ng/ul | 99     |
| 45) Dimethylphthalate   | 13.66 | 163 | 95887     | 1.631   | ng/ul | 98     |
| 50) Acenaphthene        | 14.26 | 153 | 3889336   | 79.502  | ng/ul | 99     |
| 54) Dibenzofuran        | 14.59 | 168 | 1970789   | 27.214  | ng/ul | 98     |
| 59) Fluorene            | 15.24 | 166 | 2984320   | 50.794  | ng/ul | 99     |
| 70) Phenanthrene        | 16.98 | 178 | 1790197   | 19.587  | ng/ul | 99     |
| 72) Anthracene          | 17.07 | 178 | 161784    | 1.760   | ng/ul | 97     |
| 75) Carbazole           | 17.34 | 167 | 3976097   | 51.900  | ng/ul | 100    |
| 77) Fluoranthene        | 19.00 | 202 | 710479    | 7.256   | ng/ul | 98     |
| 80) Pyrene              | 19.37 | 202 | 504457    | 4.781   | ng/ul | 99     |

(#) = .qualifier out of range (m) = manual integration (+) = signals summed