

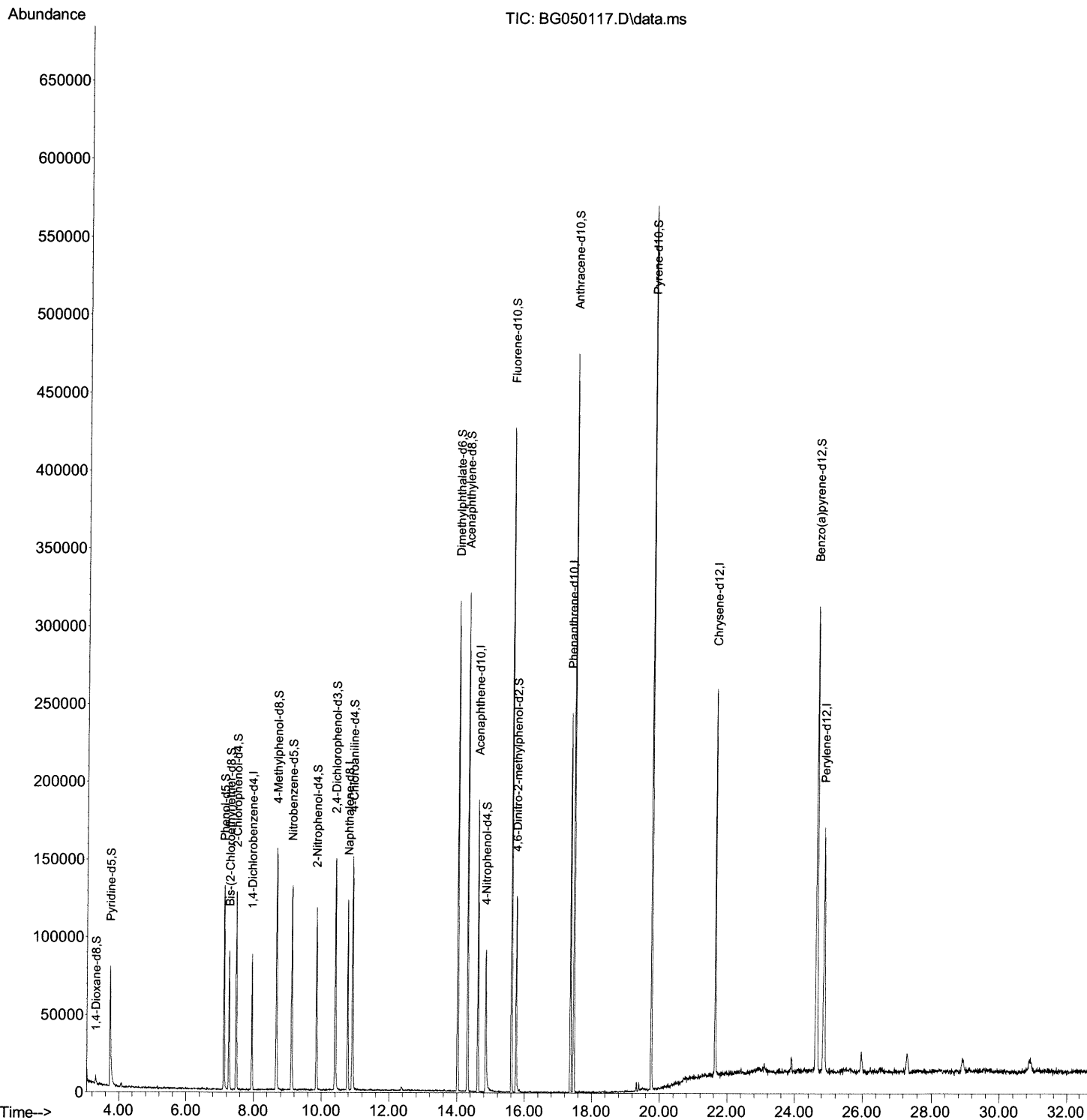
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG090221\  
Data File : BG050117.D  
Acq On : 3 Sep 2021 5:23  
Operator : CG/JU  
Sample : PB138772TB 10X  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SLEB763

Manual Integrations  
APPROVED

mohammad  
9/3/2021 3:32:29 PM

Quant Time: Sep 03 06:06:30 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG081921.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Aug 30 14:59:55 2021  
Response via : Initial Calibration



# Quantitation Report (Qedit)

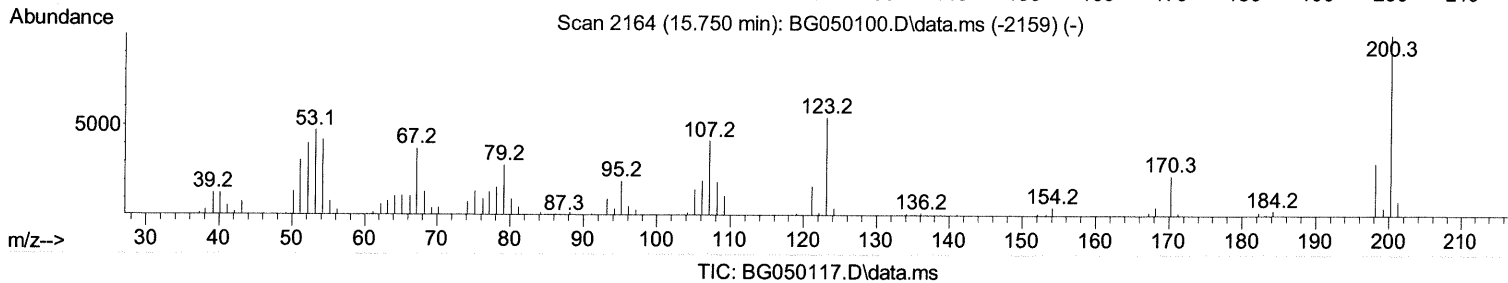
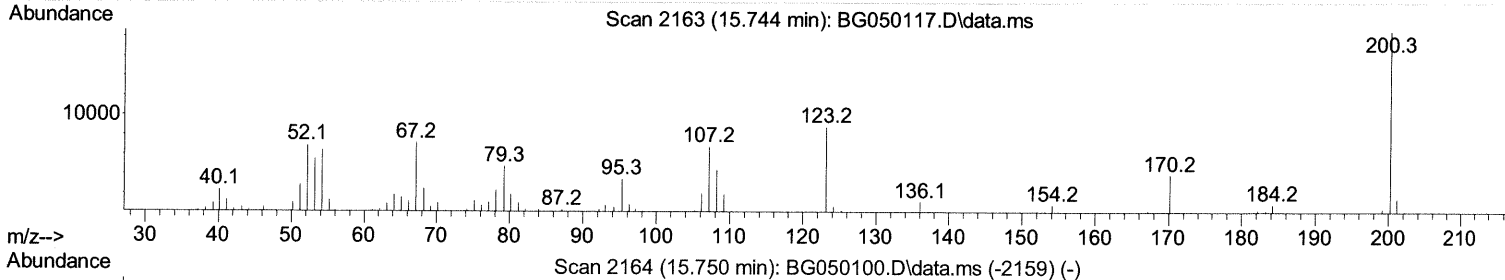
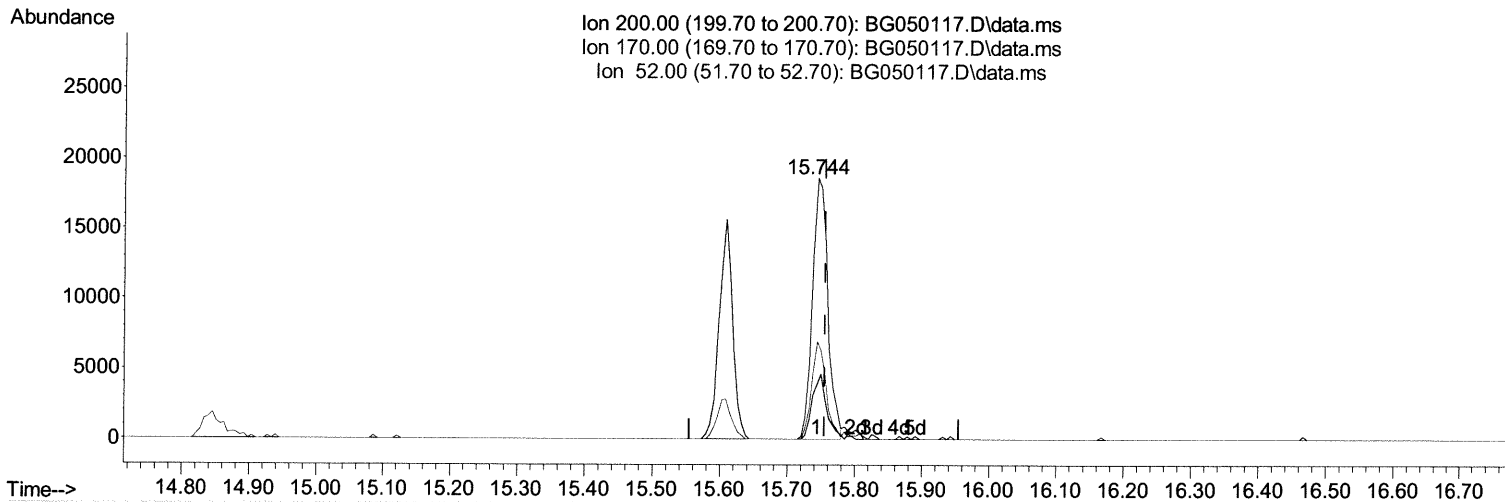
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG090221\  
 Data File : BG050117.D  
 Acq On : 3 Sep 2021 5:23  
 Operator : CG/JU  
 Sample : PB138772TB 10X  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SLEB763

Manual Integrations  
 APPROVED

mohammad  
 9/3/2021 3:32:29 PM

Quant Time: Sep 03 06:06:30 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG081921.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 30 14:59:55 2021  
 Response via : Initial Calibration



(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.744min (-0.012) 29.86 ng/ul

response 29469

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 200.00 | 100.00 | 100.00 |
| 170.00 | 22.40  | 21.22  |
| 52.00  | 45.50  | 37.31  |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

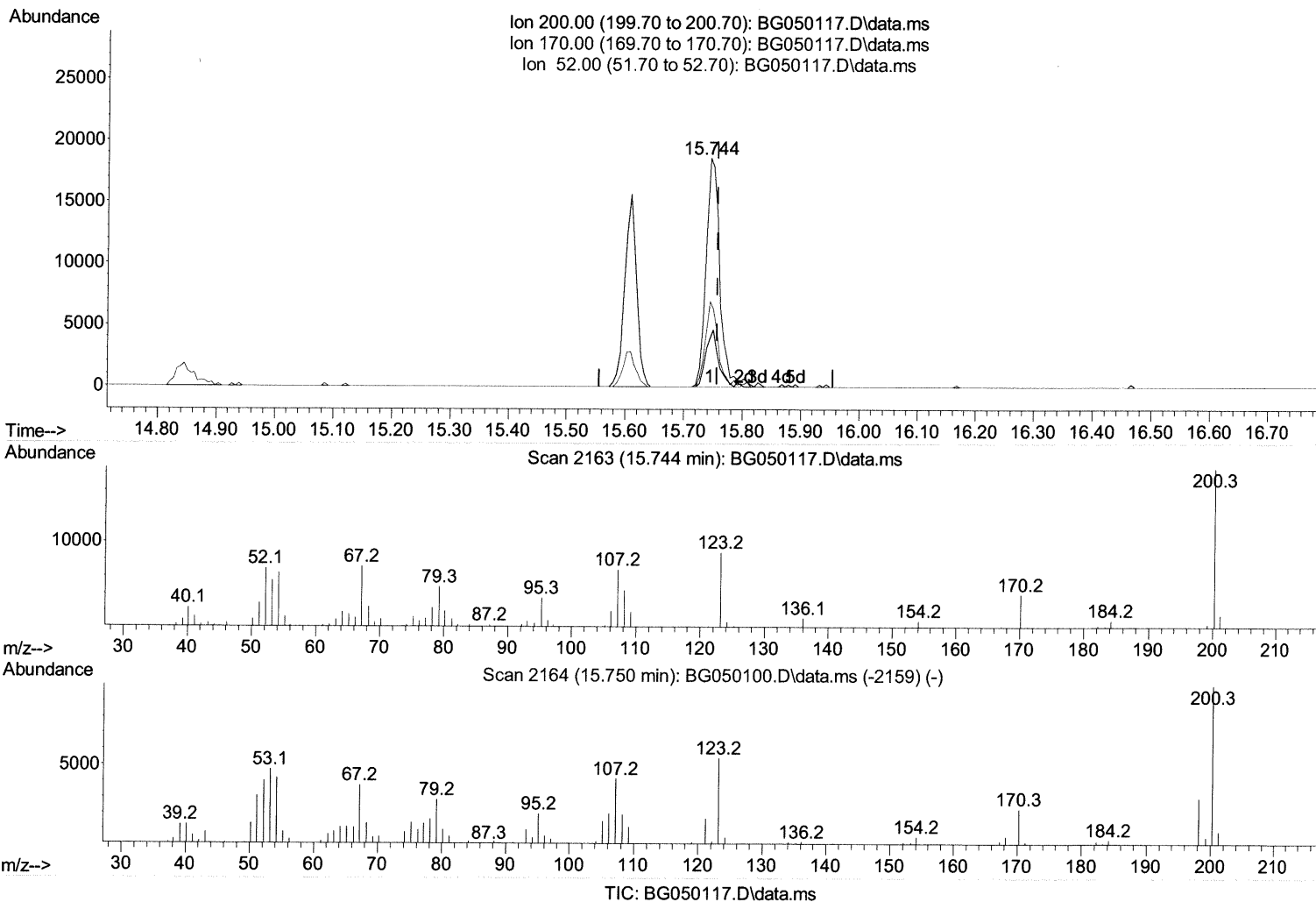
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG090221\  
 Data File : BG050117.D  
 Acq On : 3 Sep 2021 5:23  
 Operator : CG/JU  
 Sample : PB138772TB 10X  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SLEB763

Manual Integrations  
 APPROVED

mohammad  
 9/3/2021 3:32:29 PM

Quant Time: Sep 03 06:06:30 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG081921.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 30 14:59:55 2021  
 Response via : Initial Calibration



(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.744min (-0.012) 30.53 ng/ul m *54 9/8/21*

response 30127

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 200.00 | 100.00 | 100.00 |
| 170.00 | 22.40  | 21.22  |
| 52.00  | 45.50  | 37.31  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG090221\  
 Data File : BG050117.D  
 Acq On : 3 Sep 2021 5:23  
 Operator : CG/JU  
 Sample : PB138772TB 10X  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
 SLEB763

Manual Integrations  
 APPROVED

mohammad  
 9/3/2021 3:32:29 PM

Quant Time: Sep 03 06:06:30 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG081921.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 30 14:59:55 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.948  | 152  | 25496    | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.768 | 136  | 102766   | 20.000 | ng/ul | -0.01    |
| 38) Acenaphthene-d10          | 14.610 | 164  | 68514    | 20.000 | ng/ul | -0.01    |
| 64) Phenanthrene-d10          | 17.365 | 188  | 153396   | 20.000 | ng/ul | #-0.01   |
| 79) Chrysene-d12              | 21.642 | 240  | 159222   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.855 | 264  | 158012   | 20.000 | ng/ul | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.313  | 96   | 3342     | 6.759  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.736  | 84   | 54446    | 36.559 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.120  | 99   | 78733    | 36.366 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.273  | 67   | 43912    | 36.508 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.484  | 132  | 59987    | 37.030 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.671  | 113  | 60416    | 35.433 | ng/ul | -0.01    |
| 21) Nitrobenzene-d5           | 9.123  | 128  | 29884    | 37.509 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.846  | 143  | 35389    | 37.303 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.398 | 165  | 57555    | 34.014 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.909 | 131  | 80061    | 34.762 | ng/ul | -0.01    |
| 46) Dimethylphthalate-d6      | 14.011 | 166  | 202306   | 38.583 | ng/ul | -0.01    |
| 49) Acenaphthylene-d8         | 14.305 | 160  | 242337   | 39.417 | ng/ul | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.845 | 143  | 25558    | 33.690 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.609 | 176  | 168870   | 37.985 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.744 | 200  | 30127m)  | 30.528 | ng/ul | >-0.01   |
| 73) Anthracene-d10            | 17.465 | 188  | 278299   | 40.402 | ng/ul | -0.01    |
| 81) Pyrene-d10                | 19.762 | 212  | 340248   | 39.544 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.638 | 264  | 319806   | 38.149 | ng/ul | -0.01    |

Target Compounds Qvalue

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