

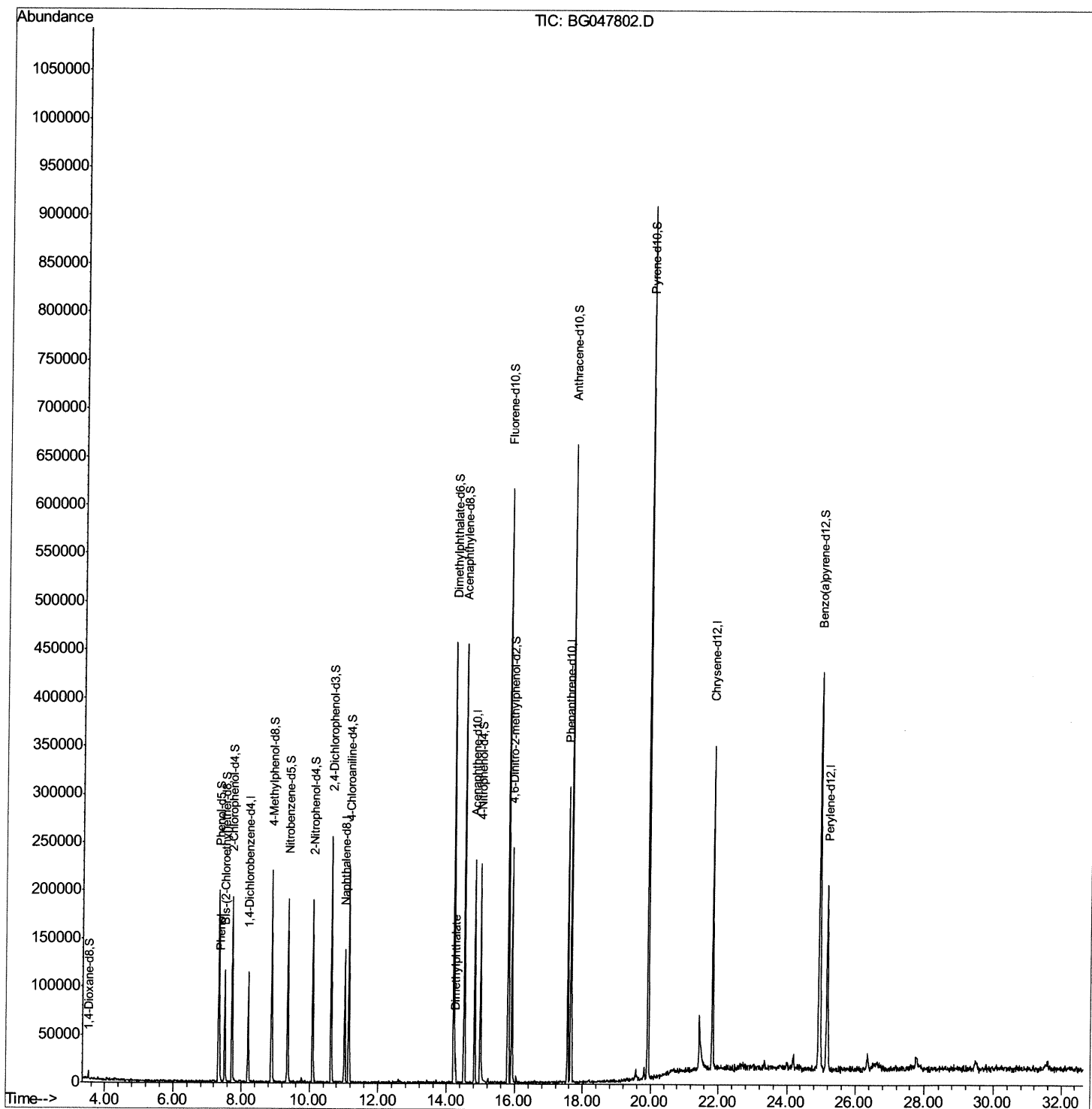
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG122320\  
Data File : BG047802.D  
Acq On : 23 Dec 2020 17:11  
Operator : CG/JU  
Sample : L5112-06  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
A54N7

Manual Integrations  
APPROVED

mohammad  
12/24/2020 10:44:09 AM

Quant Time: Dec 23 17:47:08 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG120720MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Dec 23 13:42:29 2020  
Response via : Initial Calibration



# Quantitation Report (Qedit)

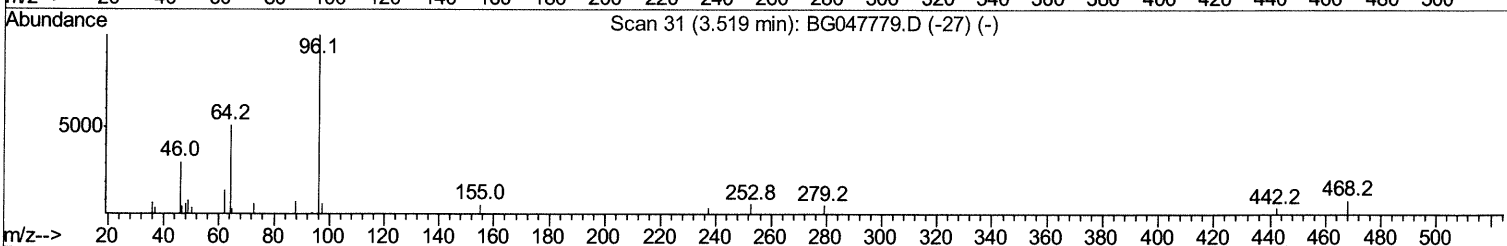
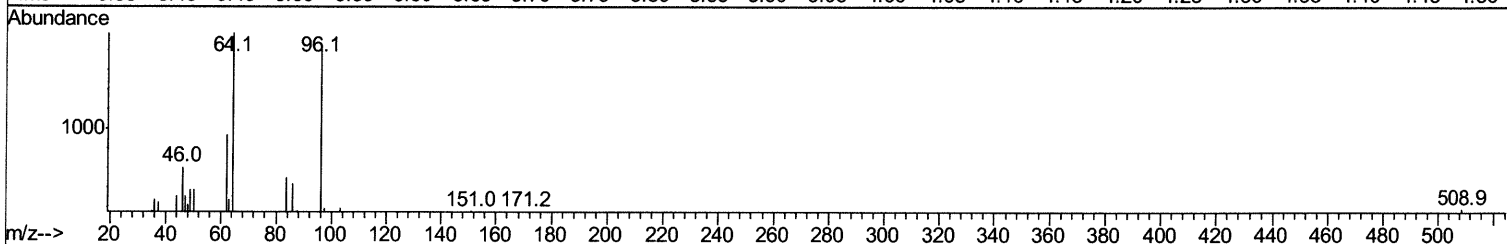
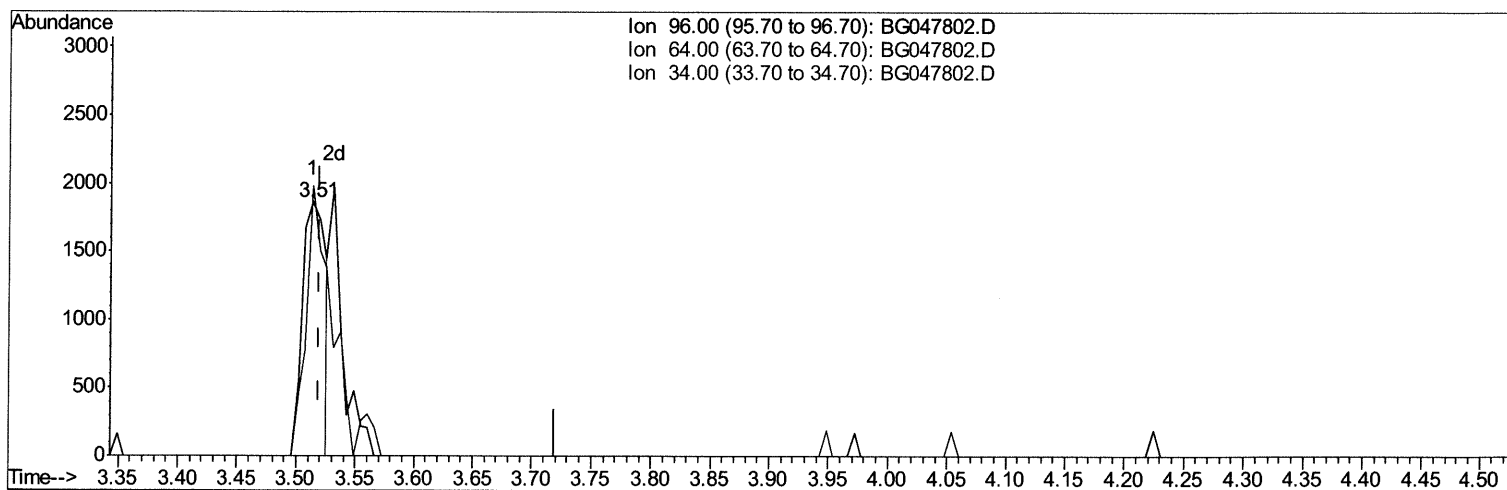
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG122320\  
 Data File : BG047802.D  
 Acq On : 23 Dec 2020 17:11  
 Operator : CG/JU  
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Manual Integrations  
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mohammad  
 12/24/2020 10:44:09 AM

Quant Time: Dec 23 17:46:36 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG120720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 23 13:42:29 2020  
 Response via : Initial Calibration



TIC: BG047802.D

(3) 1,4-Dioxane-d8 (S)

3.513min (-0.006) 5.03ng/uL

response 2562

| Ion   | Exp%  | Act%    |
|-------|-------|---------|
| 96.00 | 100   | 100     |
| 64.00 | 32.70 | 106.46# |
| 34.00 | 0.00  | 0.00    |
| 0.00  | 0.00  | 0.00    |

# Quantitation Report (Qedit)

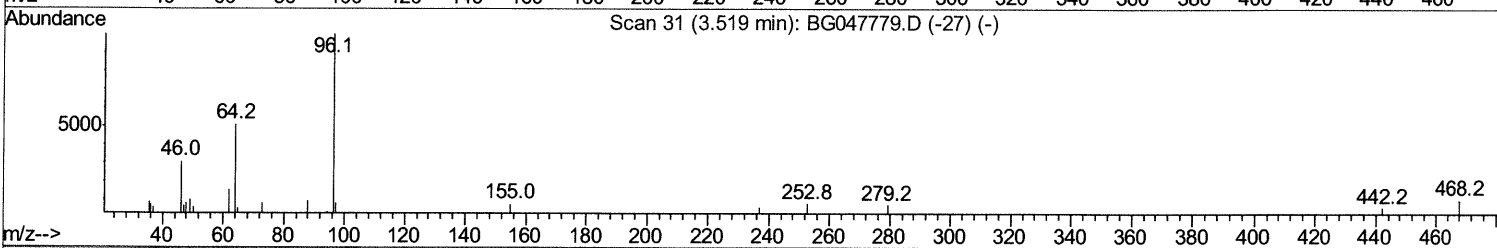
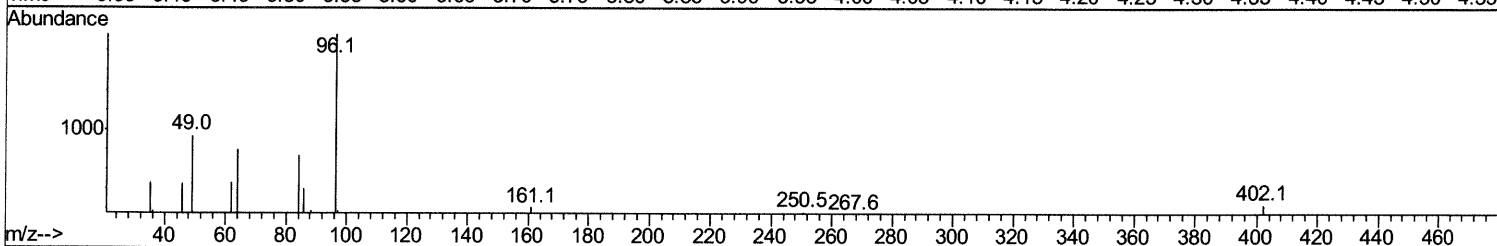
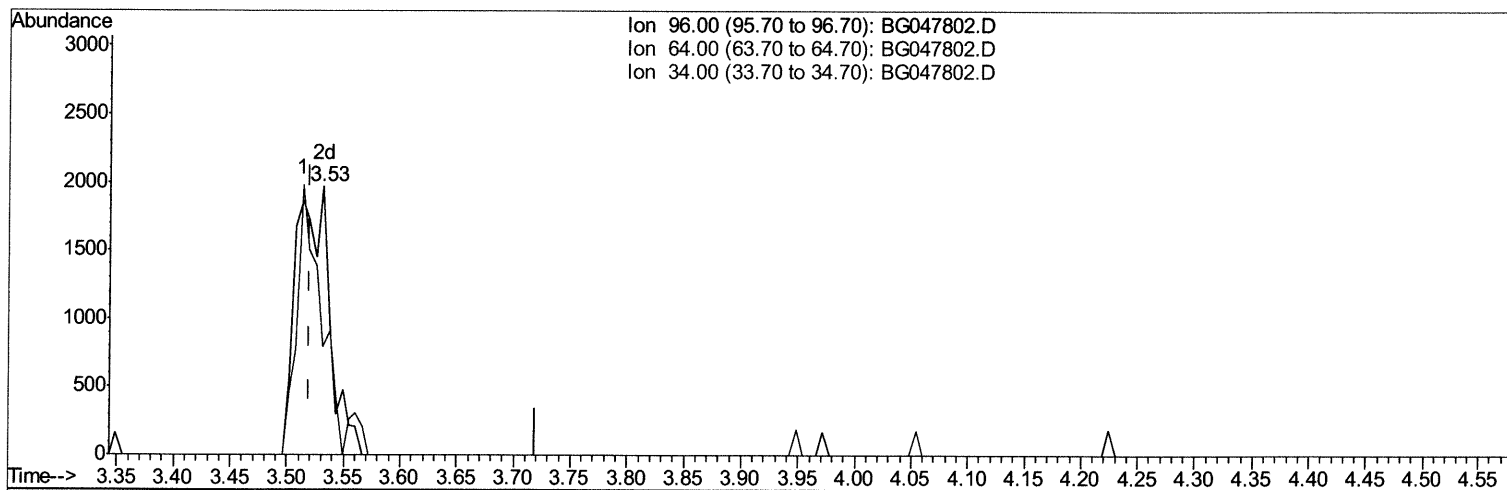
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG122320\  
 Data File : BG047802.D  
 Acq On : 23 Dec 2020 17:11  
 Operator : CG/JU  
 Sample : L5112-06  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 A54N7

Manual Integrations  
 APPROVED

mohammad  
 12/24/2020 10:44:09 AM

Quant Time: Dec 23 17:46:36 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG120720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 23 13:42:29 2020  
 Response via : Initial Calibration



TIC: BG047802.D

(3) 1,4-Dioxane-d8 (S)

3.531min (+0.012) 7.87ng/uL m 12/24/2020

response 4008

| Ion   | Exp%  | Act%   |
|-------|-------|--------|
| 96.00 | 100   | 100    |
| 64.00 | 32.70 | 40.51# |
| 34.00 | 0.00  | 0.00   |
| 0.00  | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG122320\  
 Data File : BG047802.D  
 Acq On : 23 Dec 2020 17:11  
 Operator : CG/JU  
 Sample : L5112-06  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 A54N7

Manual Integrations  
 APPROVED

mohammad  
 12/24/2020 10:44:09 AM

Quant Time: Dec 23 17:47:08 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG120720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 23 13:42:29 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.20  | 152  | 27111    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.02 | 136  | 106995   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.82 | 164  | 80932    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.56 | 188  | 195940   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.83 | 240  | 208583   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 25.17 | 264  | 210316   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |         |                  |
|--------------------------------|-------|-----|---------|-------|---------|------------------|
| 3) 1,4-Dioxane-d8              | 3.53  | 96  | 4008m > | 7.87  | ng/uL > | 0.01 12/24/20 54 |
| 5) Phenol-d5                   | 7.33  | 99  | 93256   | 40.19 | ng/ul   | 0.00             |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.51  | 67  | 50963   | 42.73 | ng/ul   | 0.00             |
| 9) 2-Chlorophenol-d4           | 7.72  | 132 | 69919   | 39.73 | ng/ul   | 0.00             |
| 13) 4-Methylphenol-d8          | 8.90  | 113 | 67254   | 36.52 | ng/ul   | 0.00             |
| 19) Nitrobenzene-d5            | 9.37  | 128 | 35077   | 40.21 | ng/ul   | 0.00             |
| 22) 2-Nitrophenol-d4           | 10.09 | 143 | 40776   | 38.78 | ng/ul   | 0.00             |
| 26) 2,4-Dichlorophenol-d3      | 10.63 | 165 | 81886   | 35.86 | ng/ul   | 0.00             |
| 29) 4-Chloroaniline-d4         | 11.15 | 131 | 95746   | 45.66 | ng/ul   | 0.00             |
| 44) Dimethylphthalate-d6       | 14.21 | 166 | 274028  | 38.98 | ng/ul   | 0.00             |
| 47) Acenaphthylene-d8          | 14.51 | 160 | 330001  | 41.10 | ng/ul   | 0.00             |
| 52) 4-Nitrophenol-d4           | 14.99 | 143 | 36602   | 35.11 | ng/ul   | 0.00             |
| 58) Fluorene-d10               | 15.80 | 176 | 243840  | 37.68 | ng/ul   | 0.00             |
| 63) 4,6-Dinitro-2-methylphenol | 15.91 | 200 | 49078   | 37.00 | ng/ul   | 0.00             |
| 71) Anthracene-d10             | 17.66 | 188 | 408492  | 41.80 | ng/ul   | 0.00             |
| 79) Pyrene-d10                 | 19.92 | 212 | 490206  | 39.76 | ng/ul   | 0.00             |
| 90) Benzo(a)pyrene-d12         | 24.95 | 264 | 492426  | 40.09 | ng/ul   | 0.00             |

## Target Compounds

|                       |       |     |       |       | Ovalue |    |
|-----------------------|-------|-----|-------|-------|--------|----|
| 6) Phenol             | 7.36  | 94  | 8578  | 3.913 | ng/ul# | 77 |
| 45) Dimethylphthalate | 14.25 | 163 | 16039 | 2.252 | ng/ul# | 90 |

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