

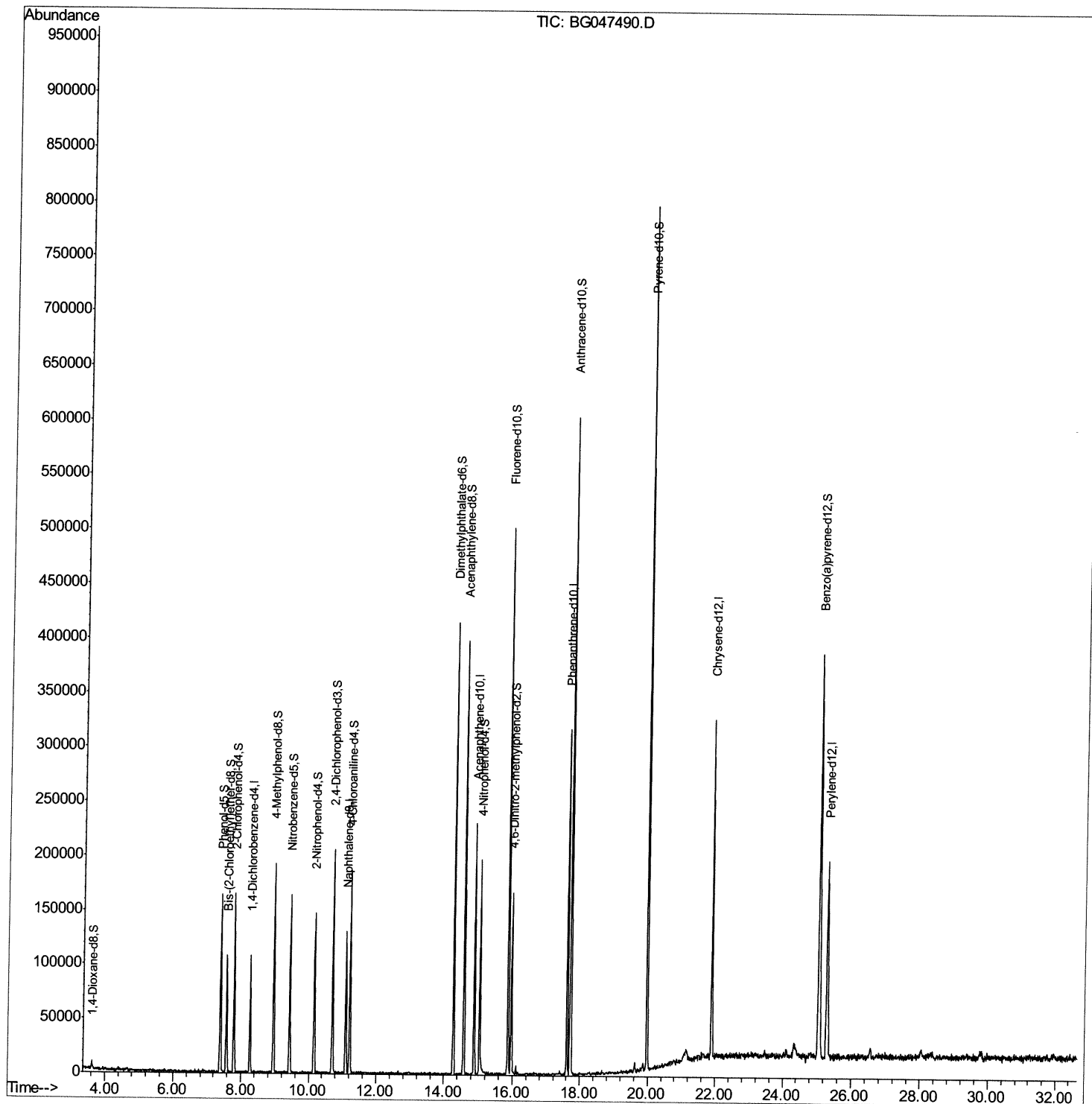
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG120120\  
Data File : BG047490.D  
Acq On : 1 Dec 2020 8:50  
Operator : CG/JU  
Sample : PB133171BL  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SBLK71

Manual Integrations  
APPROVED

mohammad  
12/2/2020 2:25:39 PM

Quant Time: Dec 01 10:16:27 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Nov 23 13:14:53 2020  
Response via : Initial Calibration



# Quantitation Report (Qedit)

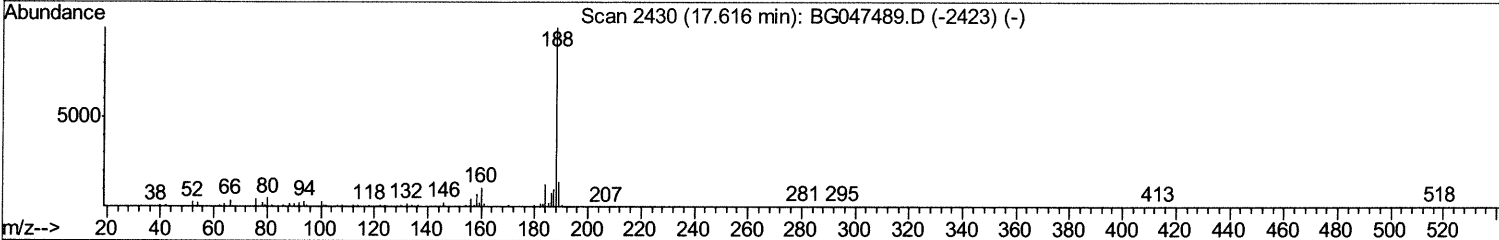
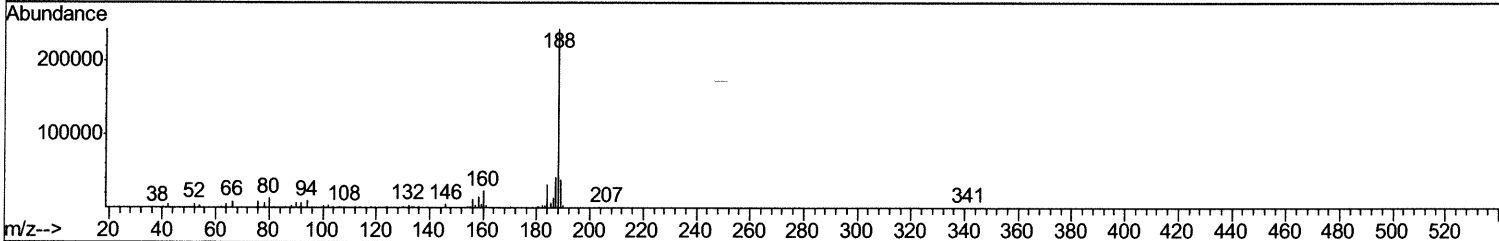
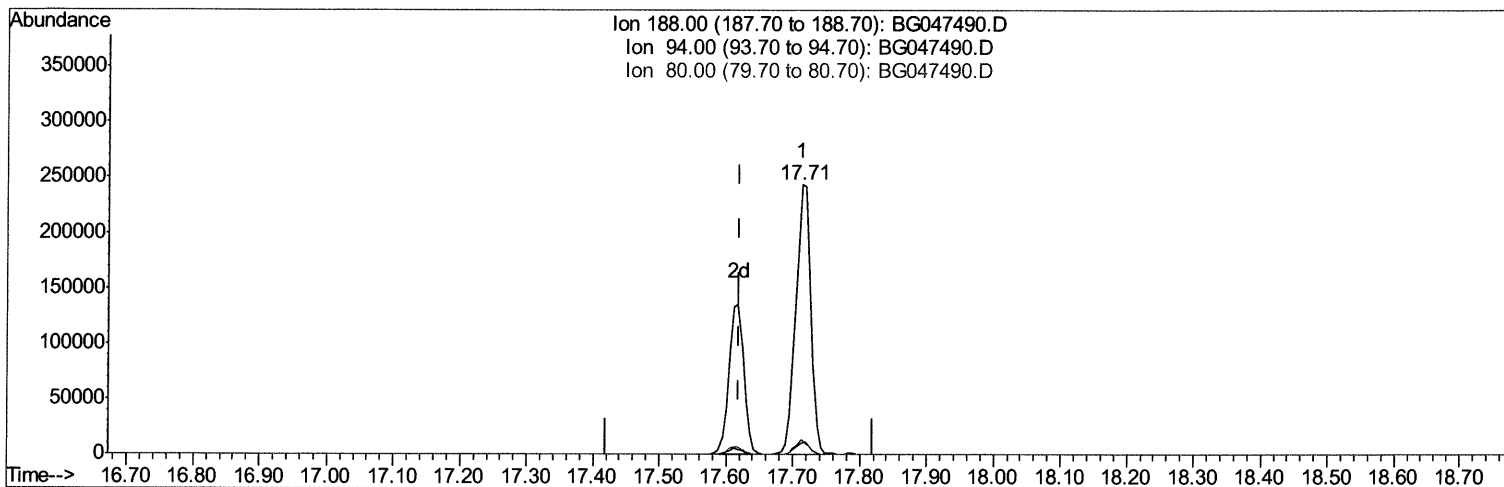
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG120120\  
 Data File : BG047490.D  
 Acq On : 1 Dec 2020 8:50  
 Operator : CG/JU  
 Sample : PB133171BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK71

Manual Integrations  
 APPROVED

mohammad  
 12/2/2020 2:25:39 PM

Quant Time: Dec 01 10:12:23 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 23 13:14:53 2020  
 Response via : Initial Calibration



TIC: BG047490.D

(62) Phenanthrene-d10 (I)

17.712min (+0.092) 20.00ng/ul

response 379522

| Ion    | Exp% | Act% |
|--------|------|------|
| 188.00 | 100  | 100  |
| 94.00  | 4.60 | 3.98 |
| 80.00  | 5.30 | 5.29 |
| 0.00   | 0.00 | 0.00 |

# Quantitation Report (Qedit)

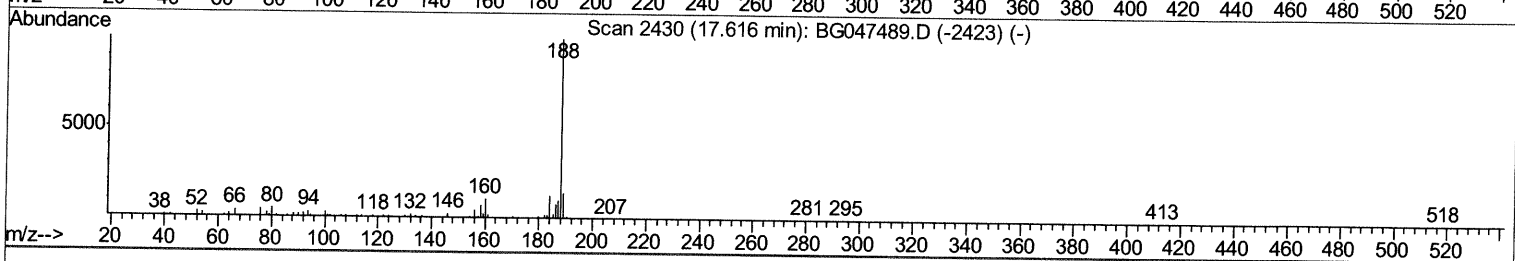
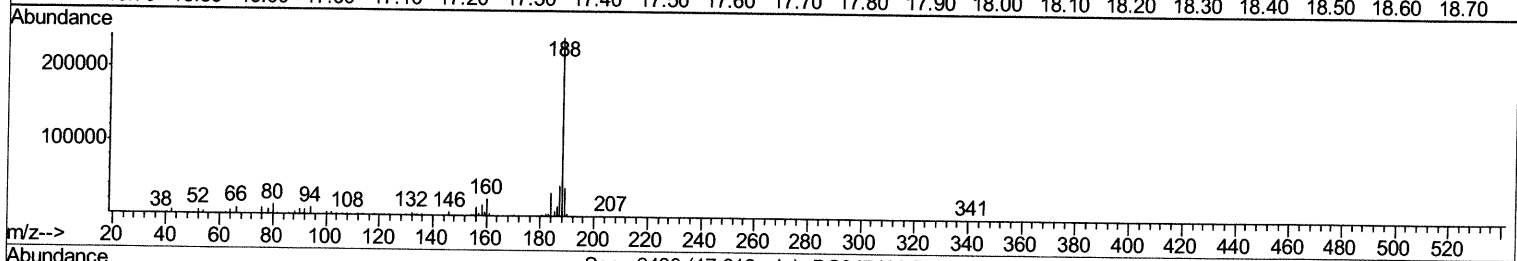
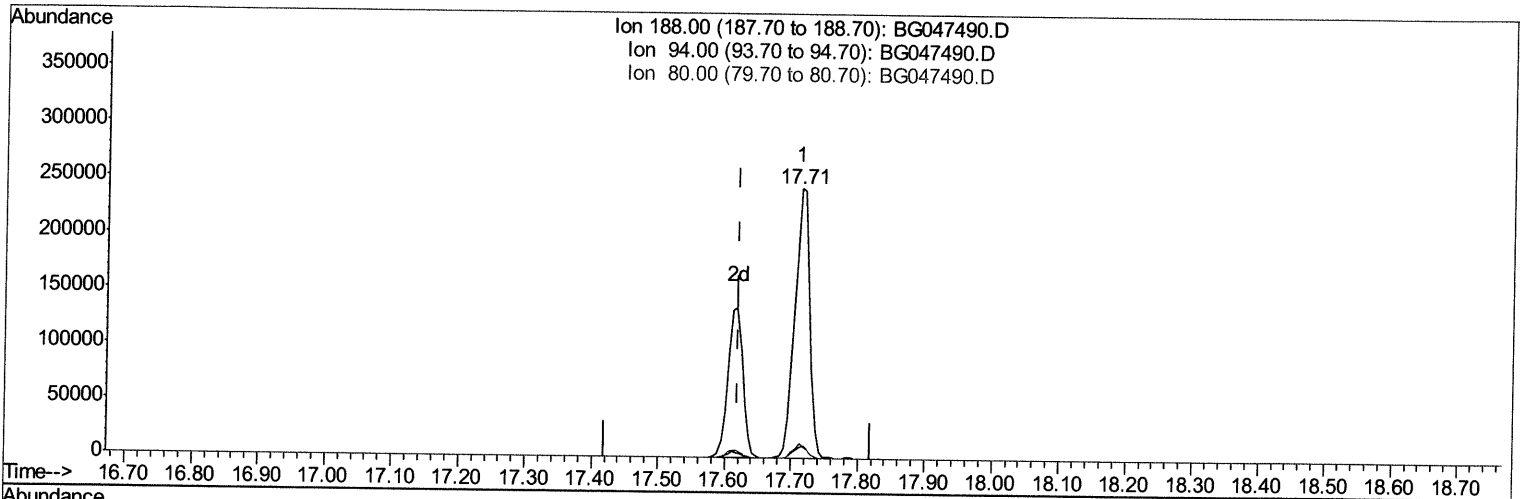
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG120120\  
 Data File : BG047490.D  
 Acq On : 1 Dec 2020 8:50  
 Operator : CG/JU  
 Sample : PB133171BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

**Instrument :**  
 BNA\_G  
**ClientSampleId :**  
 SBLK71

**Manual Integrations**  
**APPROVED**

mohammad  
 12/2/2020 2:25:39 PM

Quant Time: Dec 01 10:12:23 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 23 13:14:53 2020  
 Response via : Initial Calibration



TIC: BG047490.D

(62) Phenanthrene-d10 (I)

17.712min (+0.092) 20.00ng/ul

response 379522

| Ion    | Exp% | Act% |
|--------|------|------|
| 188.00 | 100  | 100  |
| 94.00  | 4.60 | 3.98 |
| 80.00  | 5.30 | 5.29 |
| 0.00   | 0.00 | 0.00 |

# Quantitation Report (Qedit)

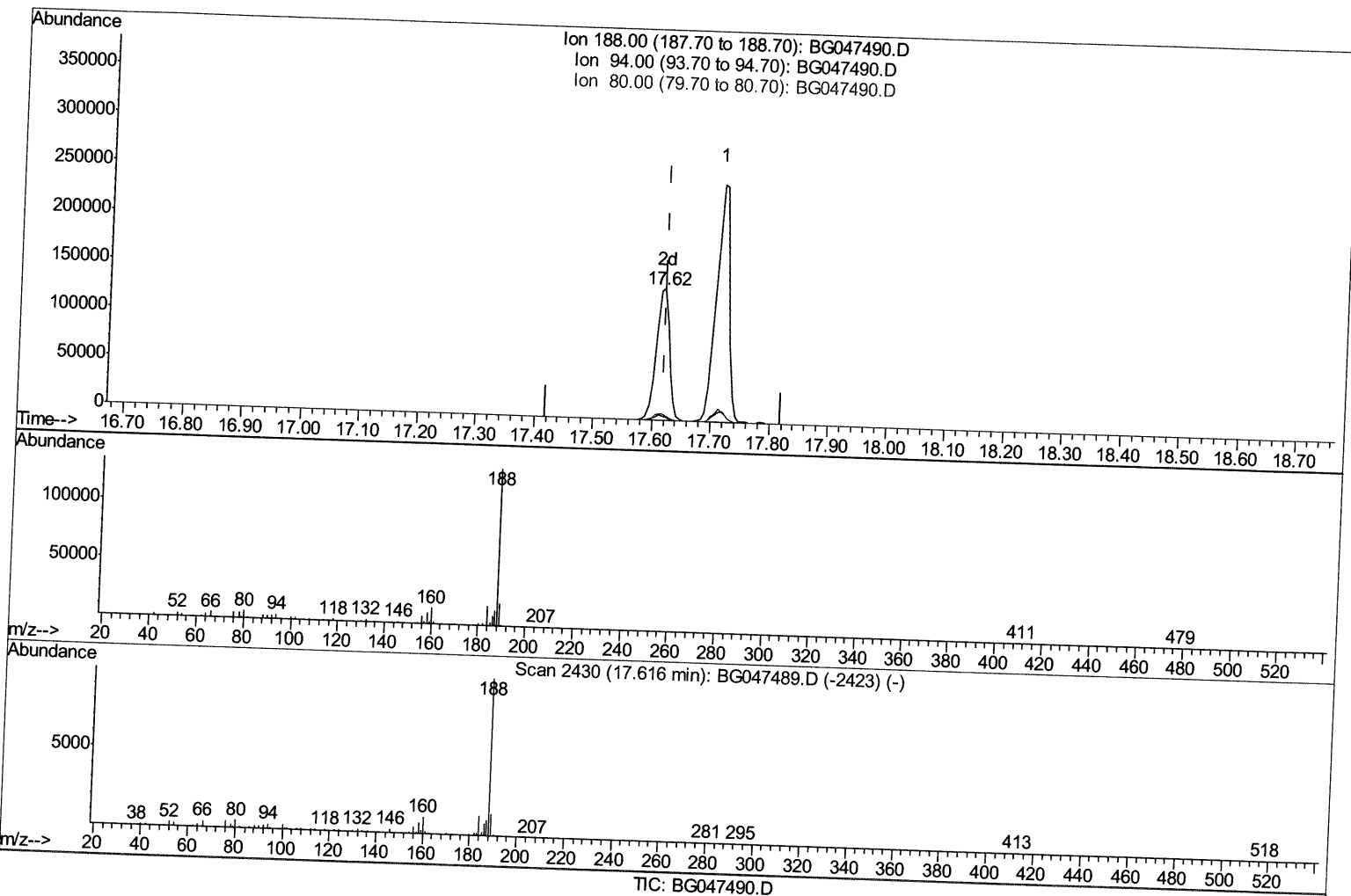
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG120120\  
 Data File : BG047490.D  
 Acq On : 1 Dec 2020 8:50  
 Operator : CG/JU  
 Sample : PB133171BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK71

Manual Integrations  
 APPROVED

mohammad  
 12/2/2020 2:25:39 PM

Quant Time: Dec 01 10:12:23 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 23 13:14:53 2020  
 Response via : Initial Calibration



(62) Phenanthrene-d10 (I)

17.618min (-0.002) 20.00ng/ul m 12/04/20 JU

response 209567

| Ion    | Exp% | Act%  |
|--------|------|-------|
| 188.00 | 100  | 100   |
| 94.00  | 4.60 | 3.10# |
| 80.00  | 5.30 | 4.63  |
| 0.00   | 0.00 | 0.00  |

## Quantitation Report (Qedit)

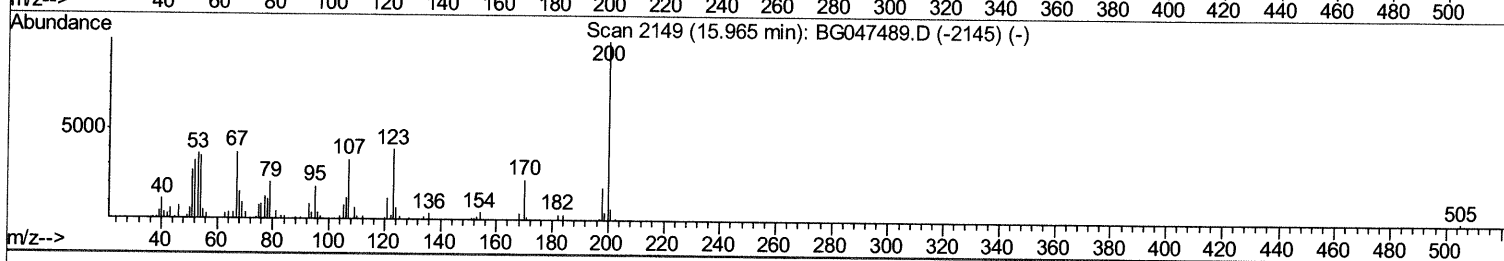
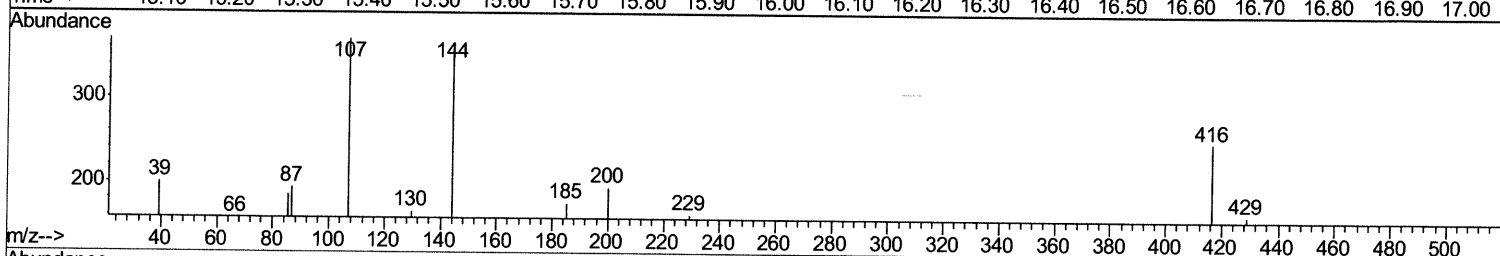
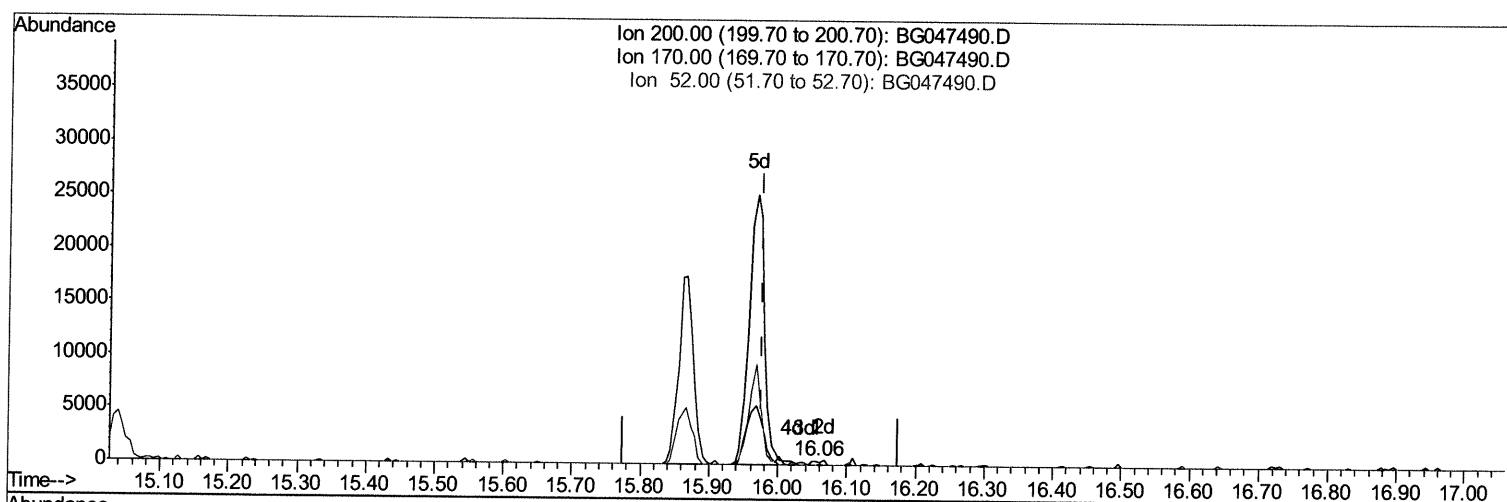
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG120120\  
Data File : BG047490.D  
Acq On : 1 Dec 2020 8:50  
Operator : CG/JU  
Sample : PB133171BL  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SBLK71

Manual Integrations  
APPROVED

mohammad  
12/2/2020 2:25:39 PM

Quant Time: Dec 01 10:15:46 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Nov 23 13:14:53 2020  
Response via : Initial Calibration



TIC: BG047490.D

(63) 4,6-Dinitro-2-methylphenol-d2 (S)

16.055min (+0.080) 0.27ng/ul

response 327

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 200.00 | 100   | 100   |
| 170.00 | 18.20 | 0.00# |
| 52.00  | 39.70 | 0.00# |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

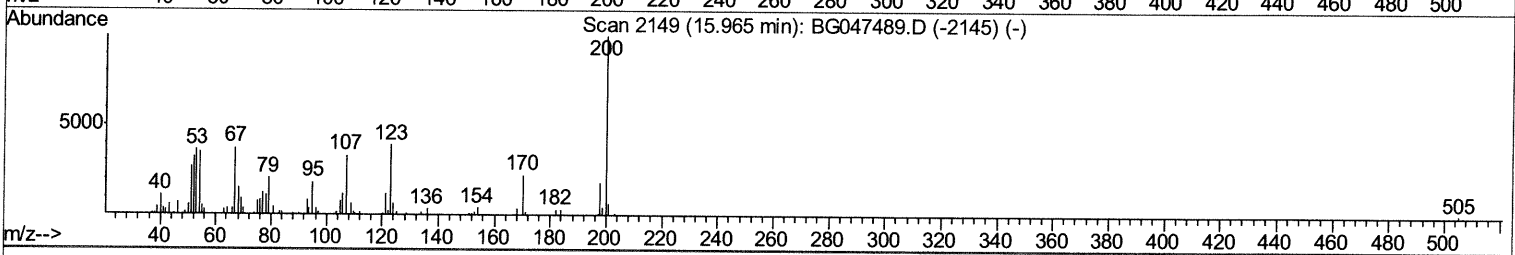
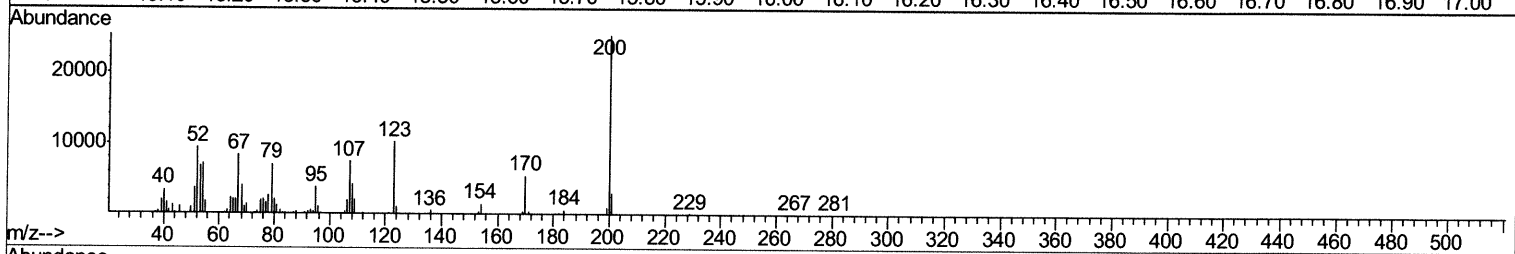
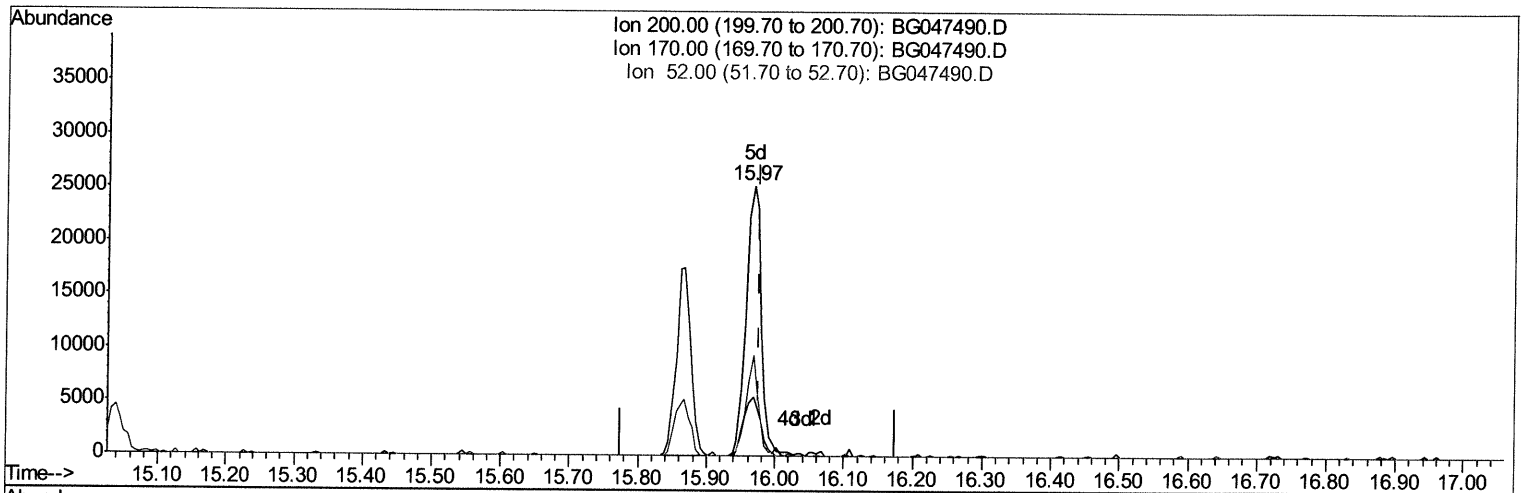
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG120120\  
 Data File : BG047490.D  
 Acq On : 1 Dec 2020 8:50  
 Operator : CG/JU  
 Sample : PB133171BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK71

Manual Integrations  
 APPROVED

mohammad  
 12/2/2020 2:25:39 PM

Quant Time: Dec 01 10:15:46 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 23 13:14:53 2020  
 Response via : Initial Calibration



TIC: BG047490.D

(63) 4,6-Dinitro-2-methylphenol-d2 (S)

15.967min (-0.008) 32.92ng/ul m 12/04/20 JU

response 39271

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 200.00 | 100   | 100    |
| 170.00 | 18.20 | 21.94# |
| 52.00  | 39.70 | 37.13  |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG120120\  
 Data File : BG047490.D  
 Acq On : 1 Dec 2020 8:50  
 Operator : CG/JU  
 Sample : PB133171BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK71

Manual Integrations  
 APPROVED

mohammad  
 12/2/2020 2:25:39 PM

Quant Time: Dec 01 10:16:27 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 23 13:14:53 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response  | Conc  | Units   | Dev(Min) |
|---------------------------|-------|------|-----------|-------|---------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.27  | 152  | 26812     | 20.00 | ng/ul   | 0.00     |
| 18) Naphthalene-d8        | 11.10 | 136  | 99906     | 20.00 | ng/ul   | 0.00     |
| 36) Acenaphthene-d10      | 14.88 | 164  | 78989     | 20.00 | ng/ul   | 0.00     |
| 62) Phenanthrene-d10      | 17.62 | 188  | 209567m > | 20.00 | ng/ul > | 0.00     |
| 78) Chrysene-d12          | 21.90 | 240  | 217554    | 20.00 | ng/ul   | 0.00     |
| 86) Perylene-d12          | 25.30 | 264  | 216680    | 20.00 | ng/ul   | 0.00     |

12/04/2020

## System Monitoring Compounds

|                                |       |     |          |       |         |       |
|--------------------------------|-------|-----|----------|-------|---------|-------|
| 3) 1,4-Dioxane-d8              | 3.61  | 96  | 3761     | 5.10  | ng/uL   | 0.00  |
| 5) Phenol-d5                   | 7.40  | 99  | 77626    | 28.87 | ng/ul   | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.58  | 67  | 46060    | 30.03 | ng/ul   | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.79  | 132 | 59227    | 32.84 | ng/ul   | 0.00  |
| 13) 4-Methylphenol-d8          | 8.96  | 113 | 60121    | 29.10 | ng/ul   | 0.00  |
| 19) Nitrobenzene-d5            | 9.43  | 128 | 29885    | 36.21 | ng/ul   | 0.00  |
| 22) 2-Nitrophenol-d4           | 10.16 | 143 | 35466    | 39.39 | ng/ul   | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 10.70 | 165 | 70177    | 34.04 | ng/ul   | 0.00  |
| 29) 4-Chloroaniline-d4         | 11.21 | 131 | 84909    | 40.73 | ng/ul   | -0.01 |
| 44) Dimethylphthalate-d6       | 14.27 | 166 | 249759   | 36.83 | ng/ul   | 0.00  |
| 47) Acenaphthylene-d8          | 14.57 | 160 | 283294   | 36.15 | ng/ul   | 0.00  |
| 52) 4-Nitrophenol-d4           | 15.04 | 143 | 35582    | 34.39 | ng/ul   | 0.00  |
| 58) Fluorene-d10               | 15.87 | 176 | 218932   | 35.25 | ng/ul   | 0.00  |
| 63) 4,6-Dinitro-2-methylphenol | 15.97 | 200 | 39271m > | 32.92 | ng/ul > | 0.00  |
| 71) Anthracene-d10             | 17.71 | 188 | 380047   | 36.56 | ng/ul   | 0.00  |
| 79) Pyrene-d10                 | 19.98 | 212 | 441316   | 34.86 | ng/ul   | 0.00  |
| 90) Benzo(a)pyrene-d12         | 25.07 | 264 | 445827   | 35.75 | ng/ul   | 0.00  |

12/04/2020

## Target Compounds

Ovalue

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