

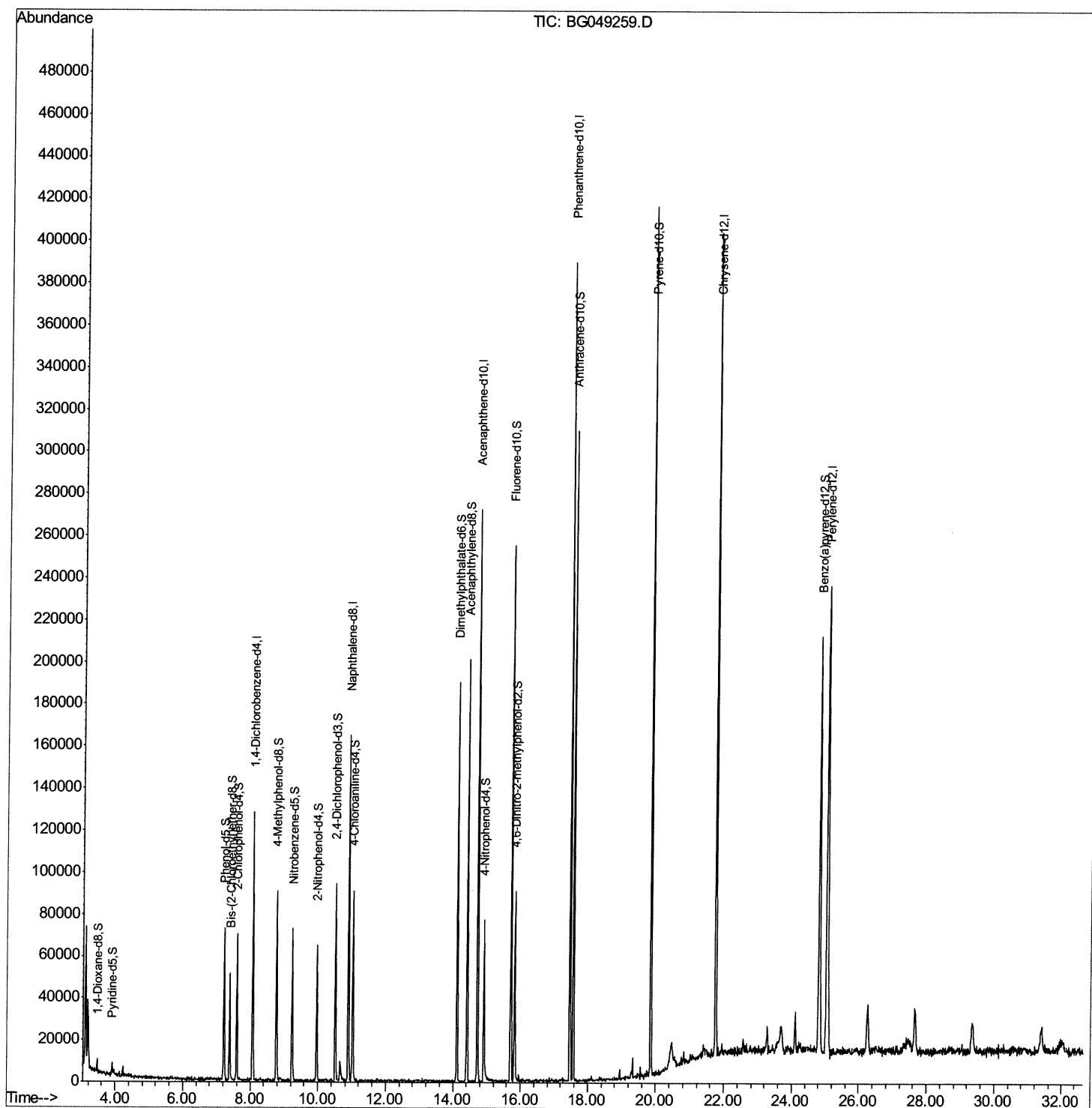
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\  
Data File : BG049259.D  
Acq On : 10 Jul 2021 10:13  
Operator : CG/JU  
Sample : M2878-02  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGDZ1

Manual Integrations  
APPROVED

mohammad  
7/12/2021 12:34:29 PM

Quant Time: Jul 11 00:51:06 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG070921.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Jul 09 03:15:38 2021  
Response via : Initial Calibration



# Quantitation Report (Qedit)

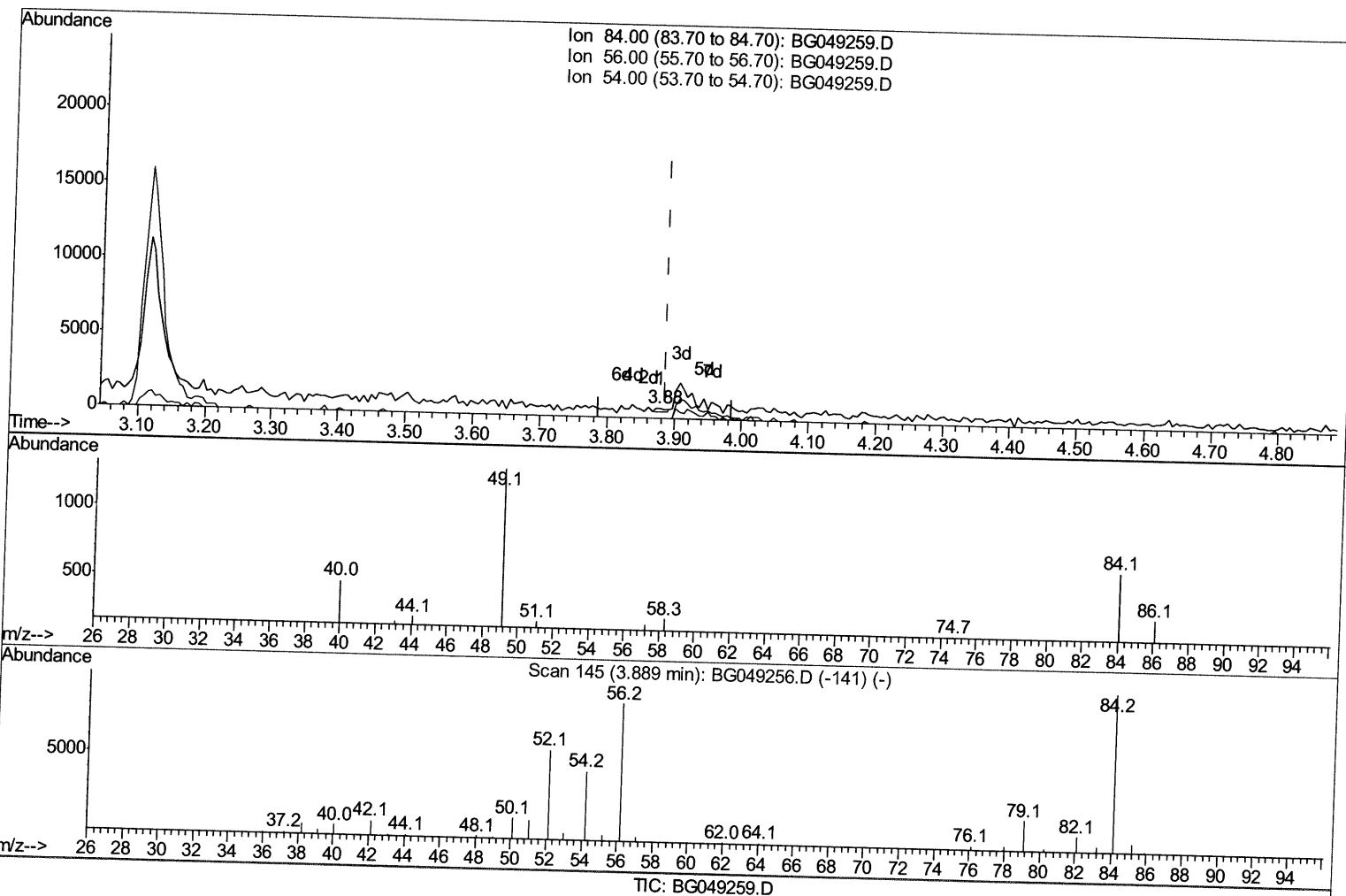
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\  
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 7/12/2021 12:34:29 PM

Quant Time: Jul 11 00:35:36 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG070921.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jul 09 03:15:38 2021  
 Response via : Initial Calibration



(4) Pyridine-d5 (S)

3.879min (-0.008) 0.14ng/ul

response 305

| Ion   | Exp%  | Act%  |
|-------|-------|-------|
| 84.00 | 100   | 100   |
| 56.00 | 76.20 | 0.00# |
| 54.00 | 42.60 | 0.00# |
| 0.00  | 0.00  | 0.00  |

# Quantitation Report (Qedit)

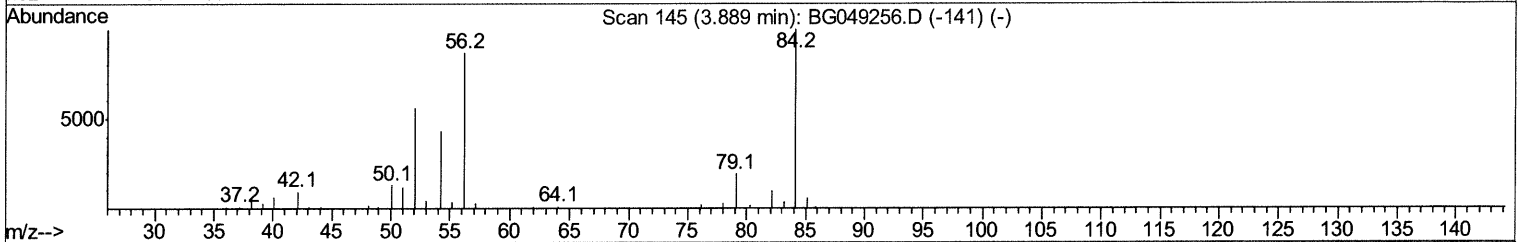
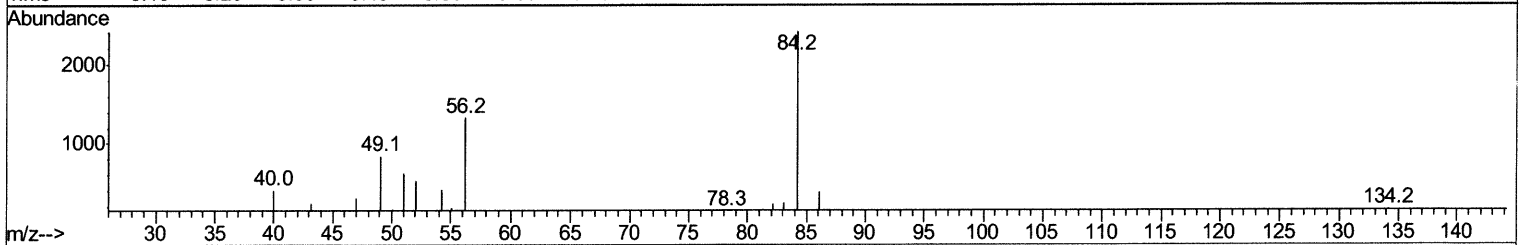
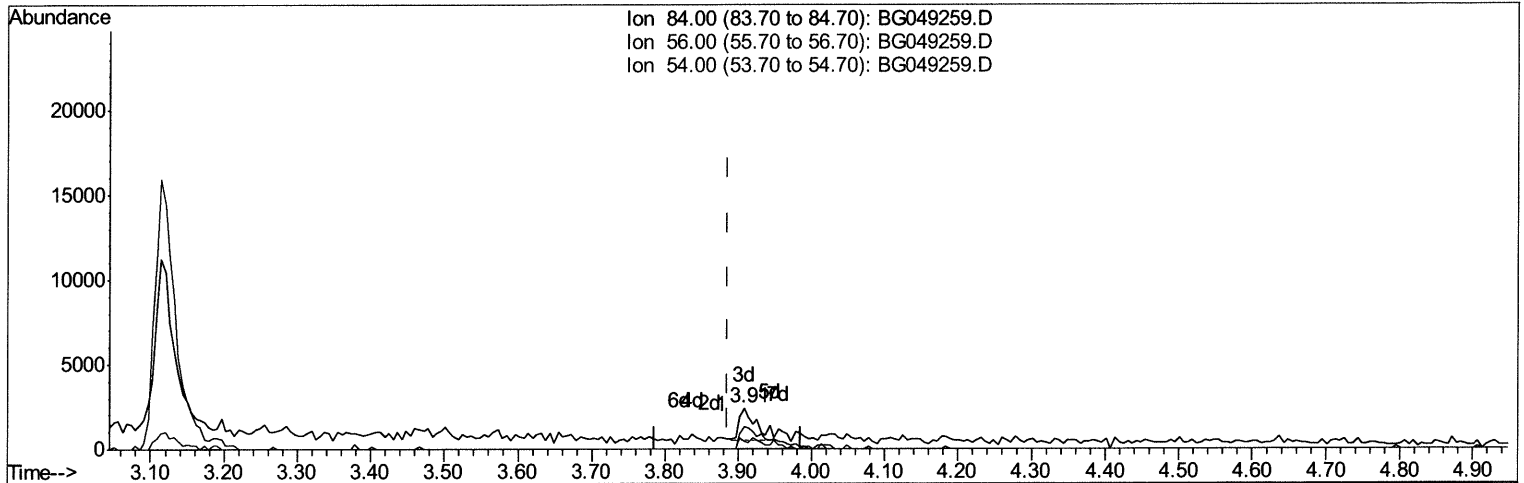
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\  
 Data File : BG049259.D  
 Acq On : 10 Jul 2021 10:13  
 Operator : CG/JU  
 Sample : M2878-02  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGDZ1

Manual Integrations  
 APPROVED

mohammad  
 7/12/2021 12:34:29 PM

Quant Time: Jul 11 00:35:36 2021  
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 Response via : Initial Calibration



TIC: BG049259.D

(4) Pyridine-d5 (S)

3.908min (+0.021) 1.46ng/ul m 07/13/21ju

response 3141

| Ion   | Exp%  | Act%   |
|-------|-------|--------|
| 84.00 | 100   | 100    |
| 56.00 | 76.20 | 54.88# |
| 54.00 | 42.60 | 17.51# |
| 0.00  | 0.00  | 0.00   |

## Quantitation Report (Qedit)

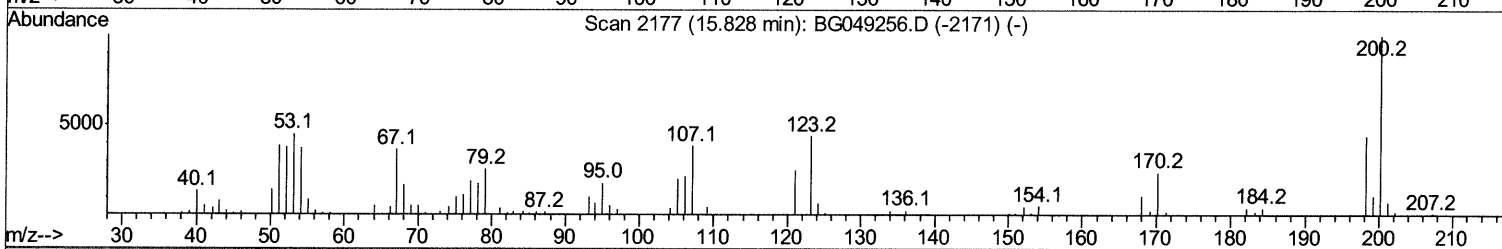
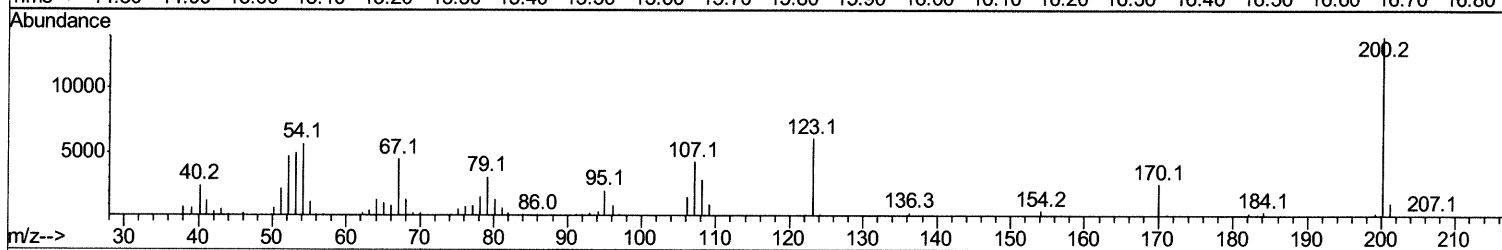
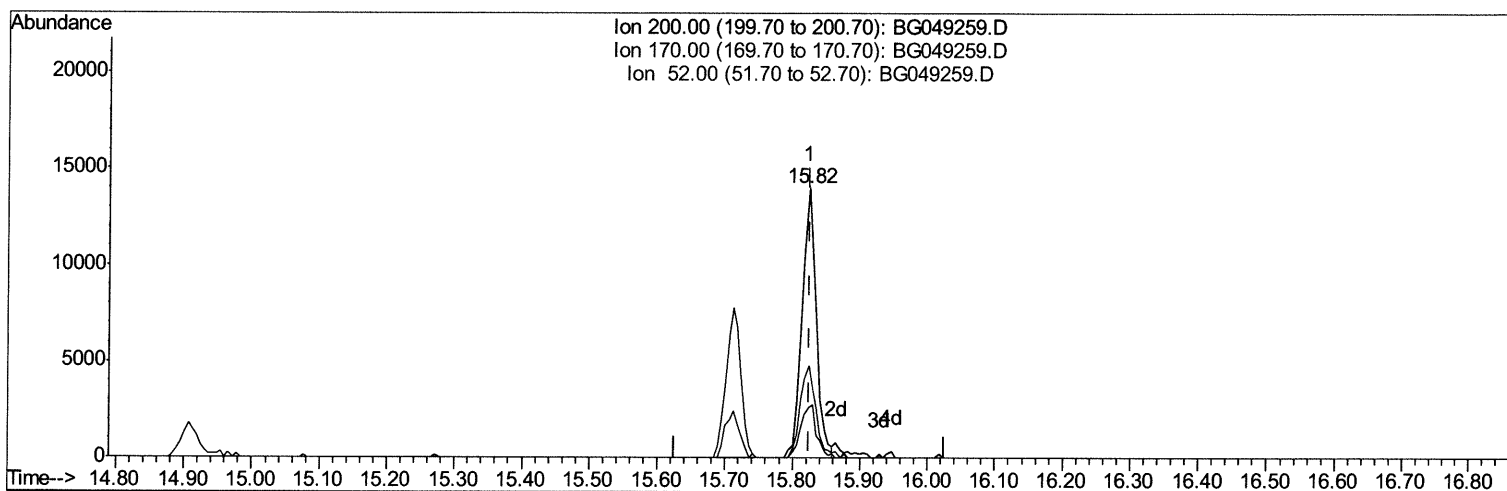
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\  
Data File : BG049259.D  
Acq On : 10 Jul 2021 10:13  
Operator : CG/JU  
Sample : M2878-02  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGDZ1

Manual Integrations  
APPROVED

mohammad  
7/12/2021 12:34:29 PM

Quant Time: Jul 11 00:50:12 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG070921.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Jul 09 03:15:38 2021  
Response via : Initial Calibration



TIC: BG049259.D

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

15.823min (-0.002) 13.29ng/ul

response 20530

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 200.00 | 100   | 100    |
| 170.00 | 23.20 | 18.45# |
| 52.00  | 47.80 | 33.94# |
| 0.00   | 0.00  | 0.00   |

# Quantitation Report (Qedit)

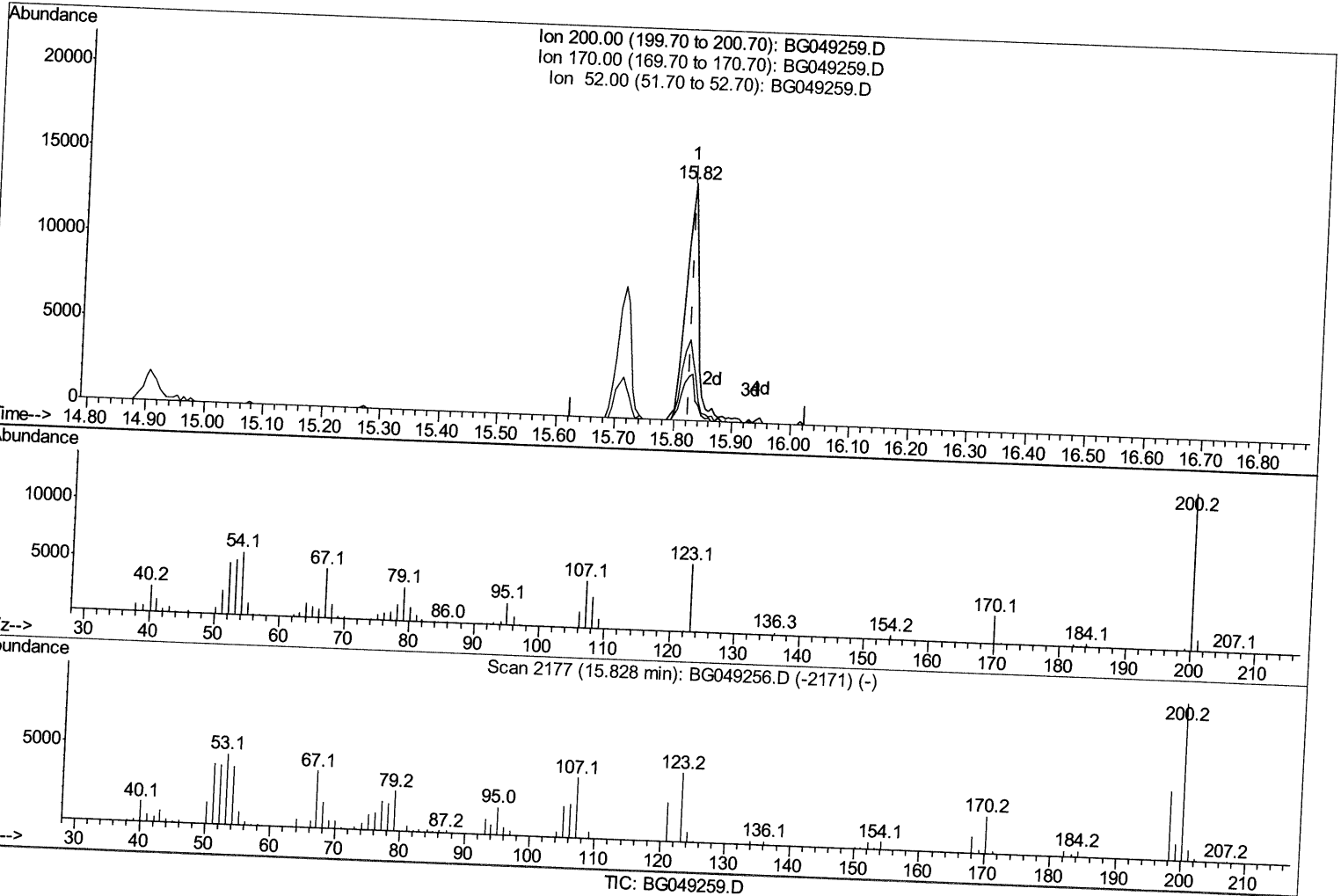
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\  
 Data File : BG049259.D  
 Acq On : 10 Jul 2021 10:13  
 Operator : CG/JU  
 Sample : M2878-02  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGDZ1

Manual Integrations  
 APPROVED

mohammad  
 7/12/2021 12:34:29 PM

Quant Time: Jul 11 00:50:12 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG070921.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jul 09 03:15:38 2021  
 Response via : Initial Calibration



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

15.823min (-0.002) 13.73ng/ul m 07/13/21 JU

response 21208

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 200.00 | 100   | 100    |
| 170.00 | 23.20 | 18.45# |
| 52.00  | 47.80 | 33.94# |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG070921\  
 Data File : BG049259.D  
 Acq On : 10 Jul 2021 10:13  
 Operator : CG/JU  
 Sample : M2878-02  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGDZ1

Manual Integrations  
 APPROVED

mohammad  
 7/12/2021 12:34:29 PM

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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.07  | 152  | 34348    | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.89 | 136  | 135581   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.72 | 164  | 100931   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.47 | 188  | 243051   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.76 | 240  | 261491   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 25.05 | 264  | 256807   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |        |      |
|--------------------------------|-------|-----|---------|-------|--------|------|
| 3) 1,4-Dioxane-d8              | 3.46  | 96  | 1693    | 2.32  | ng/uL  | 0.00 |
| 4) Pyridine-d5                 | 3.91  | 84  | 3141m>  | 1.46  | ng/ul> | 0.02 |
| 7) Phenol-d5                   | 7.22  | 99  | 37063   | 12.40 | ng/ul  | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.39  | 67  | 24816   | 14.32 | ng/ul  | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.60  | 132 | 28404   | 12.86 | ng/ul  | 0.00 |
| 15) 4-Methylphenol-d8          | 8.77  | 113 | 32348   | 13.51 | ng/ul  | 0.00 |
| 21) Nitrobenzene-d5            | 9.24  | 128 | 15269   | 14.68 | ng/ul  | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.97  | 143 | 19170   | 16.05 | ng/ul  | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.51 | 165 | 34379   | 14.75 | ng/ul  | 0.00 |
| 31) 4-Chloroaniline-d4         | 11.03 | 131 | 44600   | 14.69 | ng/ul  | 0.00 |
| 46) Dimethylphthalate-d6       | 14.12 | 166 | 126596  | 16.31 | ng/ul  | 0.00 |
| 49) Acenaphthylene-d8          | 14.41 | 160 | 140399  | 15.42 | ng/ul  | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.91 | 143 | 18799   | 14.63 | ng/ul  | 0.00 |
| 60) Fluorene-d10               | 15.71 | 176 | 107018  | 16.28 | ng/ul  | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.82 | 200 | 21208m> | 13.73 | ng/ul> | 0.00 |
| 73) Anthracene-d10             | 17.57 | 188 | 187911  | 16.98 | ng/ul  | 0.00 |
| 81) Pyrene-d10                 | 19.85 | 212 | 227817  | 17.00 | ng/ul  | 0.00 |
| 92) Benzo(a)pyrene-d12         | 24.82 | 264 | 227956  | 17.08 | ng/ul  | 0.00 |

## Target Compounds

Ovalue

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