

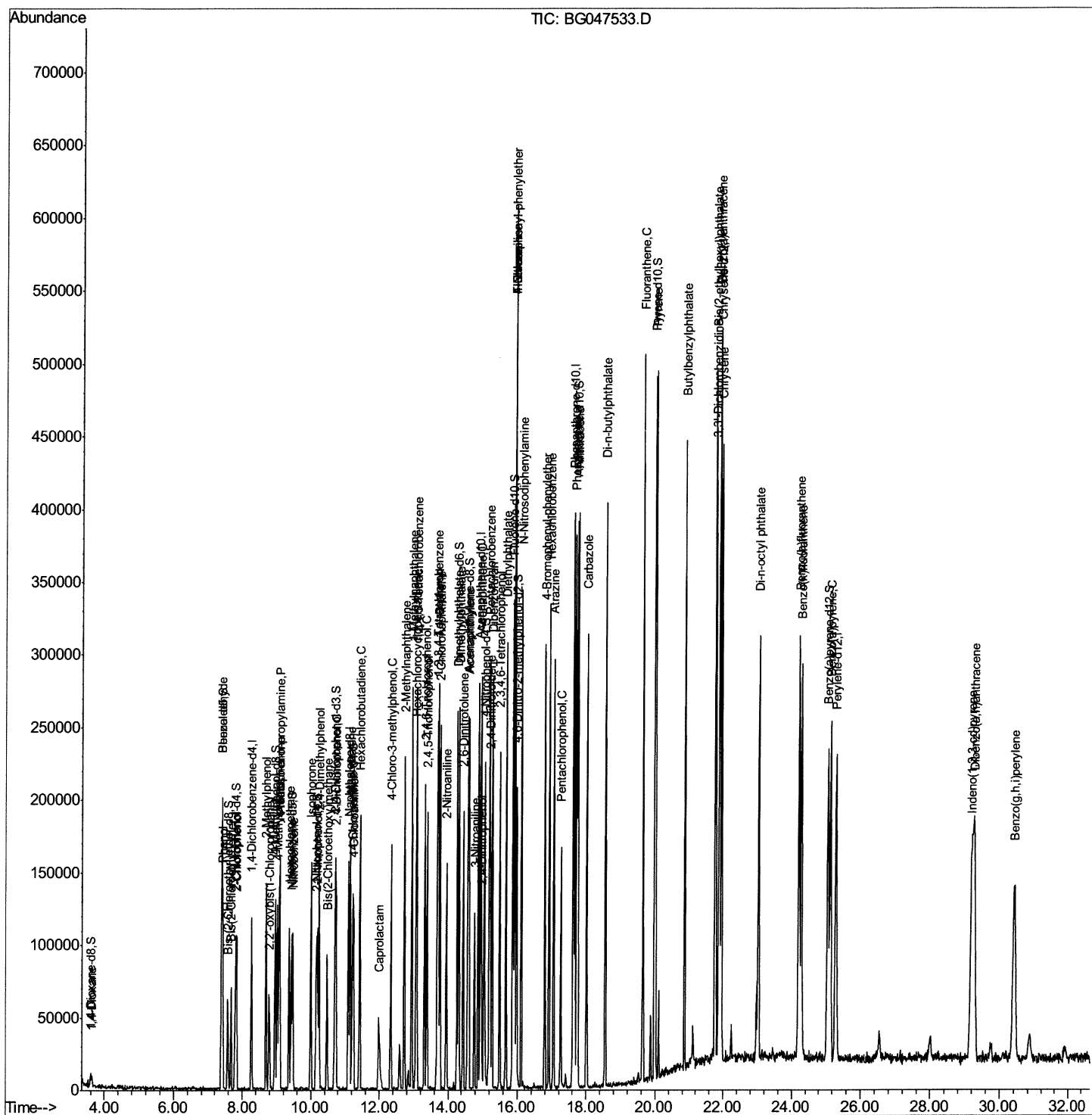
```
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
Data File : BG047533.D  
Acq On    : 2 Dec 2020 16:04  
Operator  : CG/JU  
Sample    : SSTDCCC020  
Misc      :  
ALS Vial  : 46 Sample Multiplier: 1
```

**Instrument :**  
BNA\_G  
**LabSampleId :**  
SSTD02069

## Manual Integrations APPROVED

mohammad  
12/3/2020 3:35:14 PM

Quant Time: Dec 02 17:02:46 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



# Quantitation Report (Qedit)

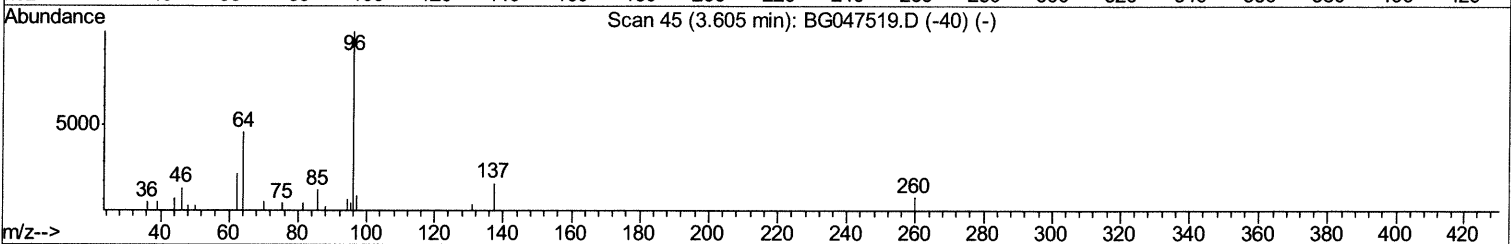
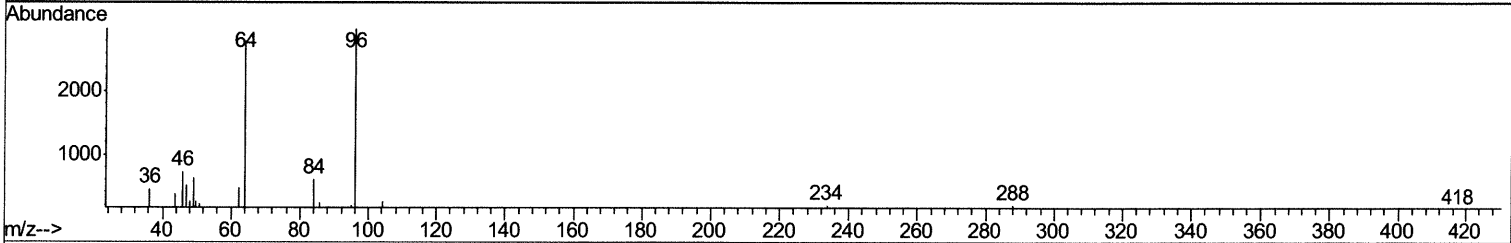
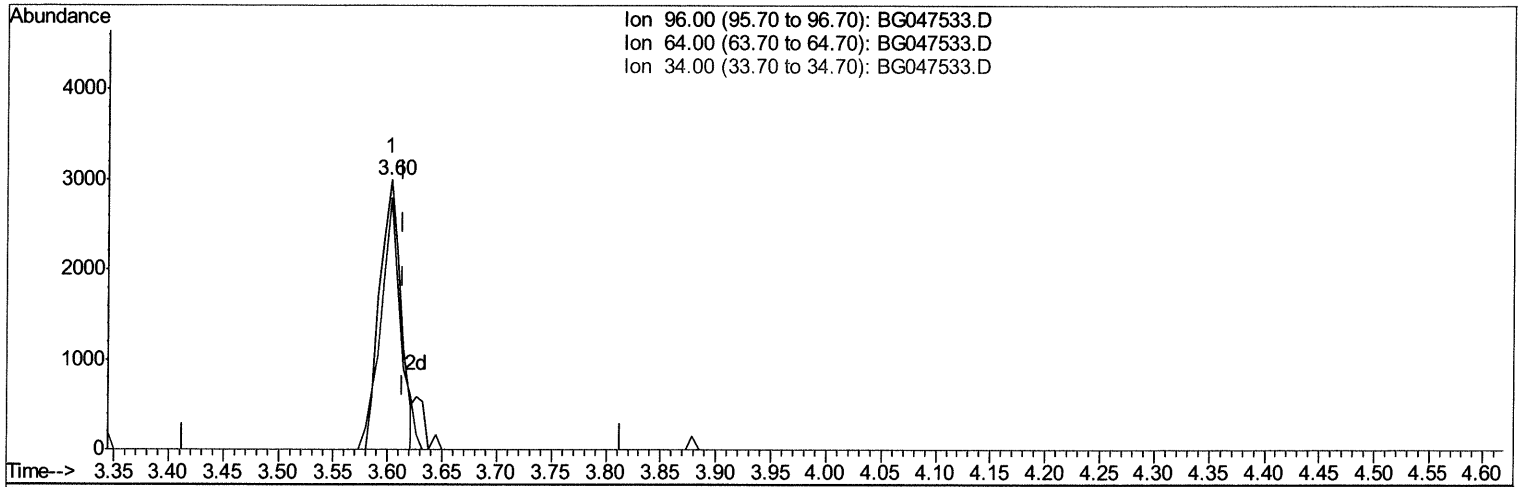
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
 Data File : BG047533.D  
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 Operator : CG/JU  
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 Misc :  
 ALS Vial : 46 Sample Multiplier: 1

**Instrument :**  
 BNA\_G  
**LabSampleId :**  
 SSTD02069

**Manual Integrations**  
**APPROVED**

mohammad  
 12/3/2020 3:35:14 PM

Quant Time: Dec 02 16:59:37 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: BG047533.D

(3) 1,4-Dioxane-d8 (S)  
 3.603min (-0.010) 4.93ng/uL  
 response 4024

| Ion   | Exp%  | Act%  |
|-------|-------|-------|
| 96.00 | 100   | 100   |
| 64.00 | 88.60 | 93.29 |
| 34.00 | 0.00  | 0.00  |
| 0.00  | 0.00  | 0.00  |

## Quantitation Report (Qedit)

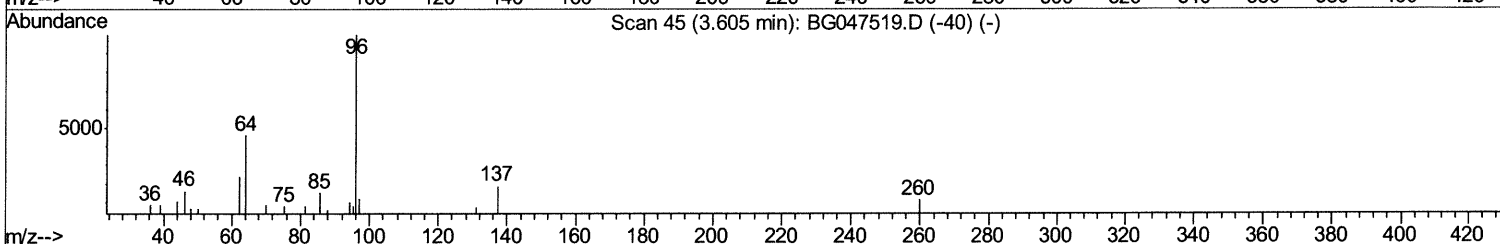
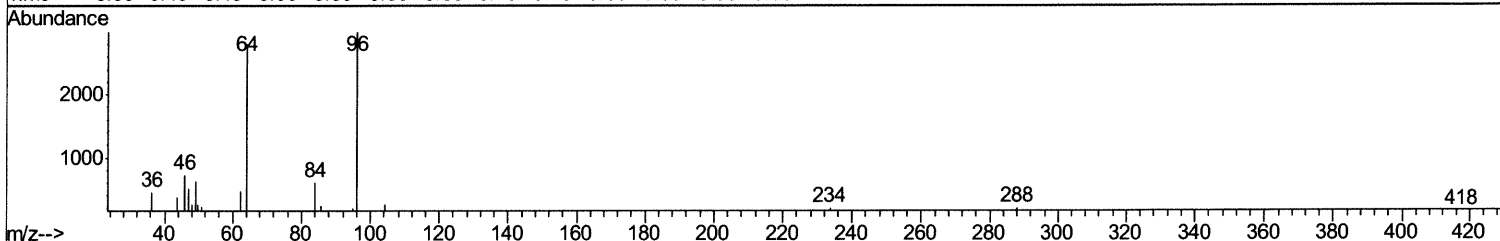
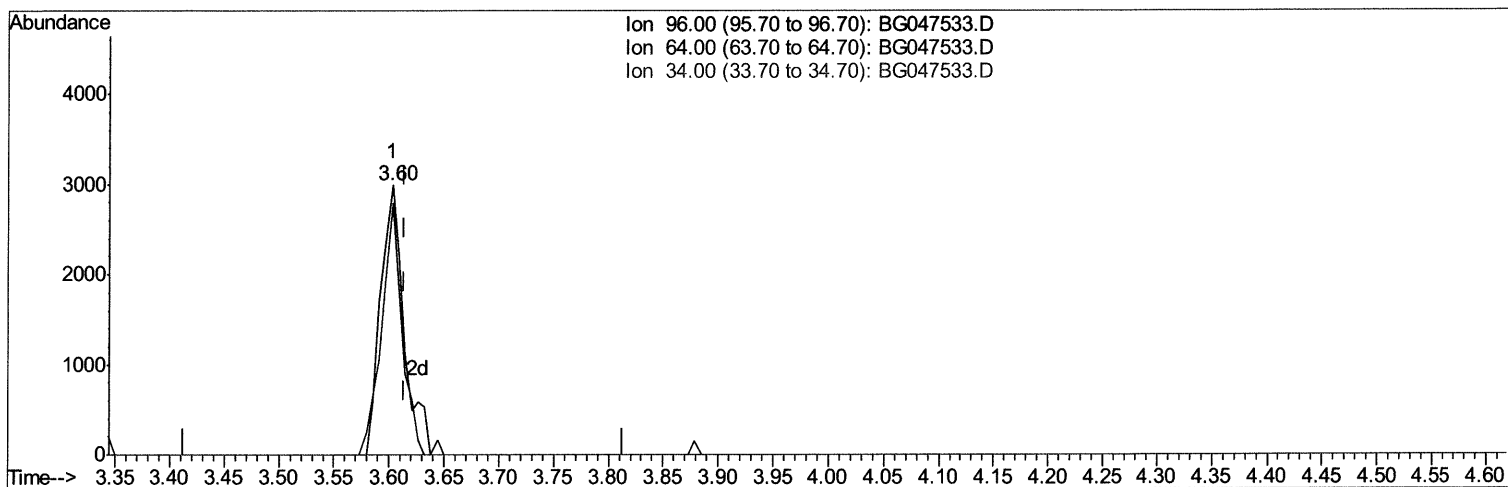
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
Data File : BG047533.D  
Acq On : 2 Dec 2020 16:04  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02069

Manual Integrations  
APPROVED

mohammad  
12/3/2020 3:35:14 PM

Quant Time: Dec 02 16:59:37 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: BG047533.D

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

3.603min (-0.010) 5.41ng/uL m 12/04/20 JU

response 4416

| Ion   | Exp%  | Act%  |
|-------|-------|-------|
| 96.00 | 100   | 100   |
| 64.00 | 88.60 | 93.29 |
| 34.00 | 0.00  | 0.00  |
| 0.00  | 0.00  | 0.00  |

# Quantitation Report (Qedit)

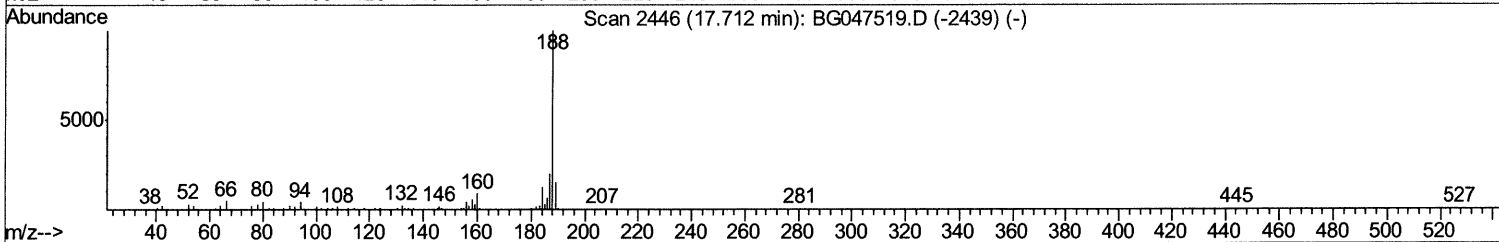
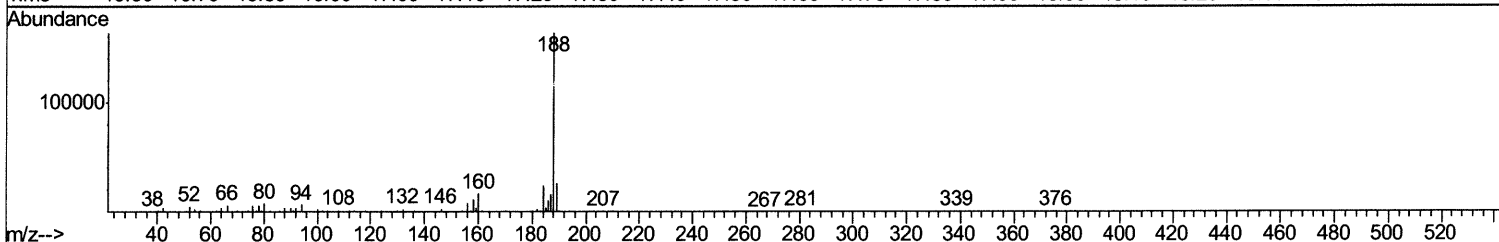
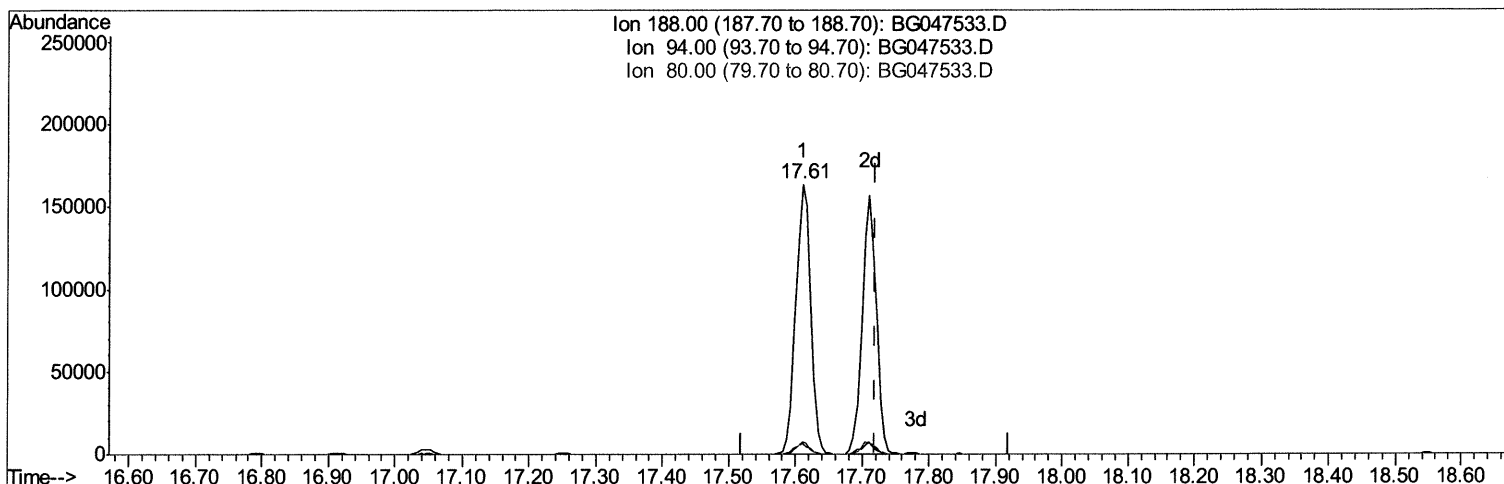
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
 Data File : BG047533.D  
 Acq On : 2 Dec 2020 16:04  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 46 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02069

Manual Integrations  
 APPROVED

mohammad  
 12/3/2020 3:35:14 PM

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 Quant Title : SVOA CALIBRATION  
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 Response via : Initial Calibration



TIC: BG047533.D

(71) Anthracene-d10 (S)

17.610min (-0.110) 20.16ng/ul

response 254498

| Ion    | Exp% | Act% |
|--------|------|------|
| 188.00 | 100  | 100  |
| 94.00  | 4.50 | 4.00 |
| 80.00  | 5.10 | 4.61 |
| 0.00   | 0.00 | 0.00 |

## Quantitation Report (Qedit)

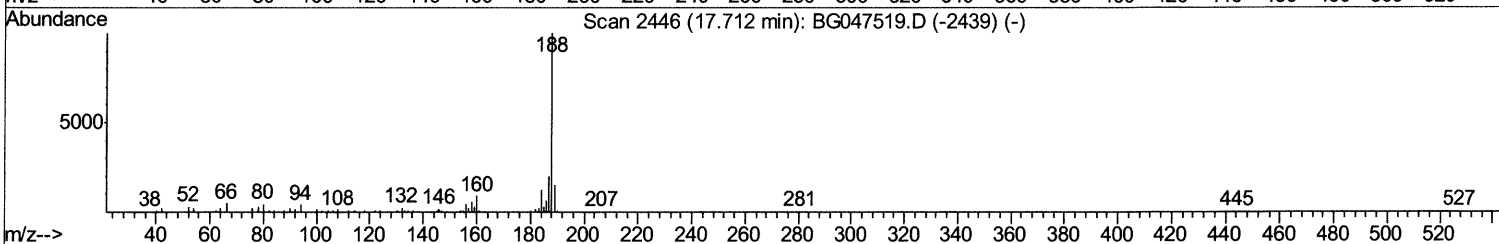
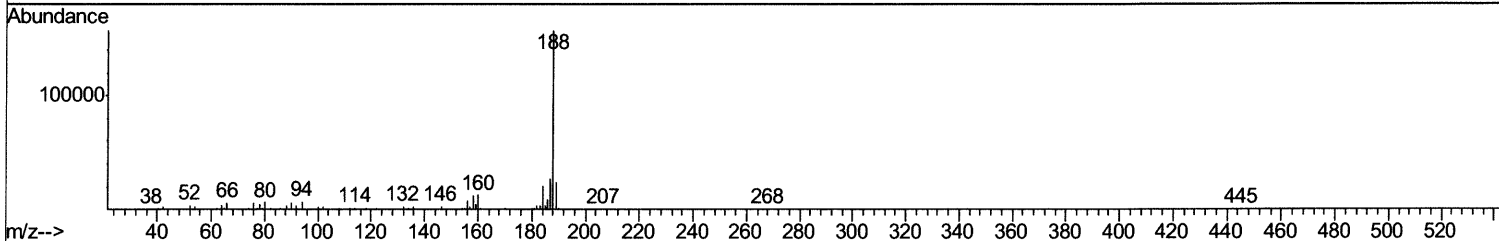
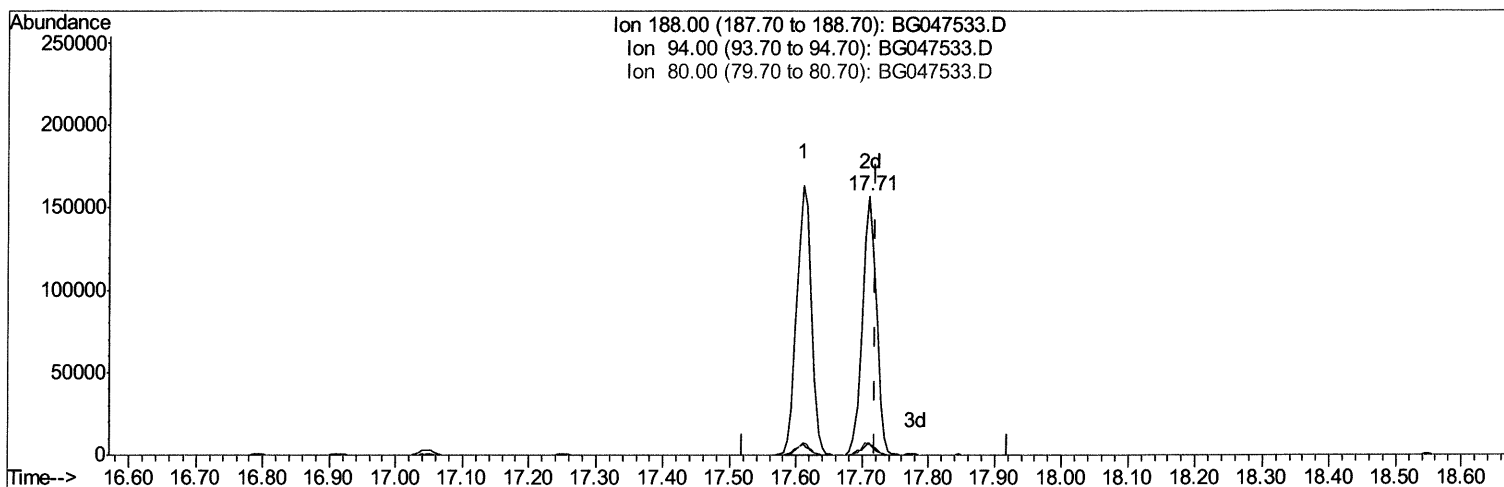
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
Data File : BG047533.D  
Acq On : 2 Dec 2020 16:04  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02069

Manual Integrations  
APPROVED

mohammad  
12/3/2020 3:35:14 PM

Quant Time: Dec 02 16:59:37 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: BG047533.D

(71) Anthracene-d10 (S)

17.710min (-0.010) 18.46ng/ul m 12/04/20 JU

response 233027

| Ion    | Exp% | Act%  |
|--------|------|-------|
| 188.00 | 100  | 100   |
| 94.00  | 4.50 | 4.53  |
| 80.00  | 5.10 | 4.07# |
| 0.00   | 0.00 | 0.00  |

## Quantitation Report (Qedit)

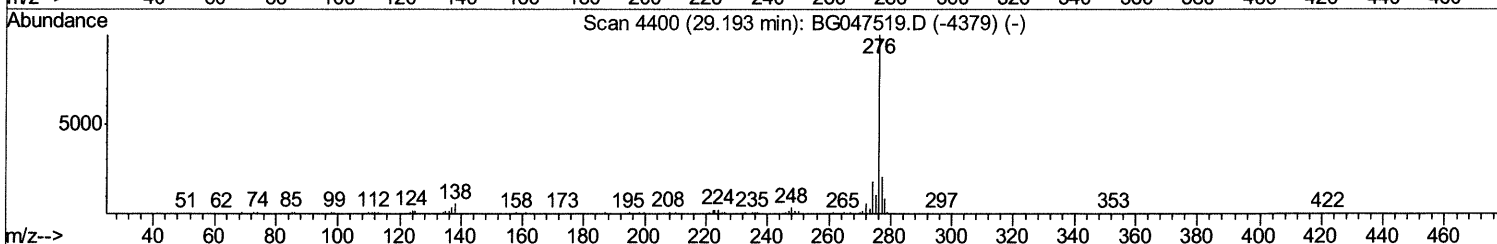
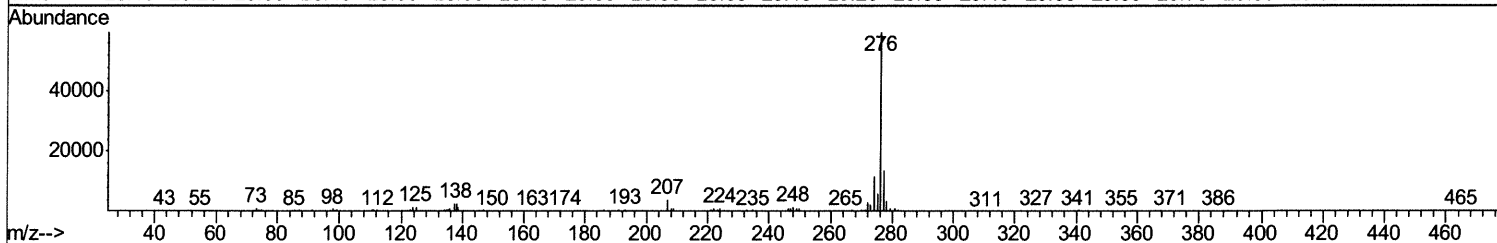
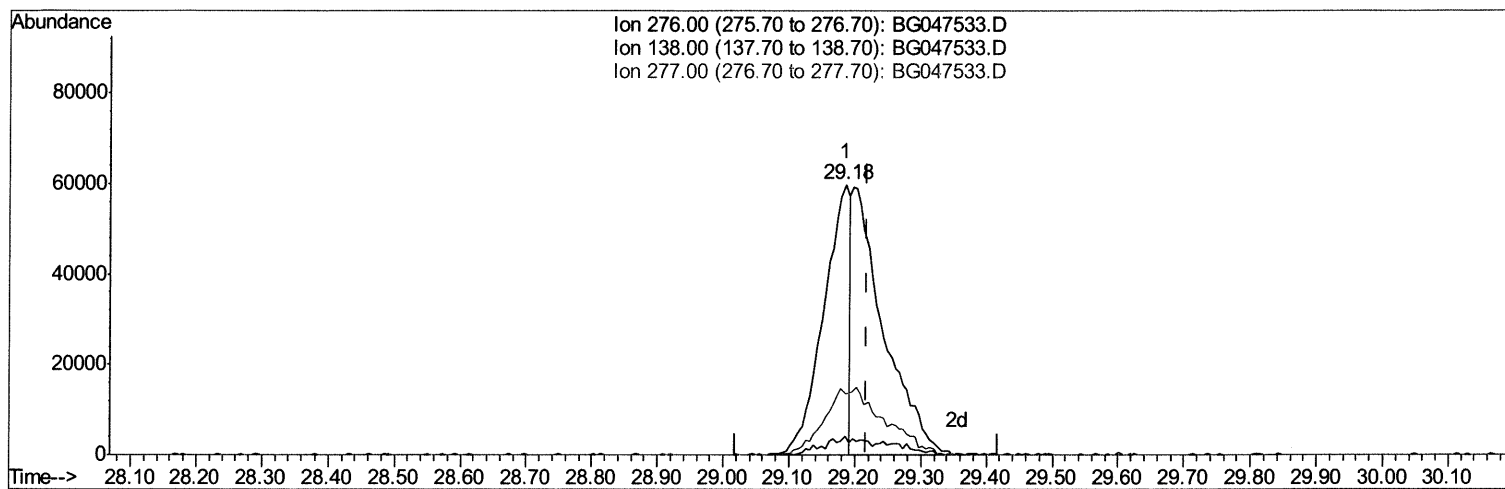
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
Data File : BG047533.D  
Acq On : 2 Dec 2020 16:04  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02069

Manual Integrations  
APPROVED

mohammad  
12/3/2020 3:35:14 PM

Quant Time: Dec 02 16:59:37 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: BG047533.D

(92) Indeno(1,2,3-cd)pyrene

29.185min (-0.034) 8.55ng/ul

response 163811

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 6.30  | 6.84  |
| 277.00 | 20.60 | 22.82 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

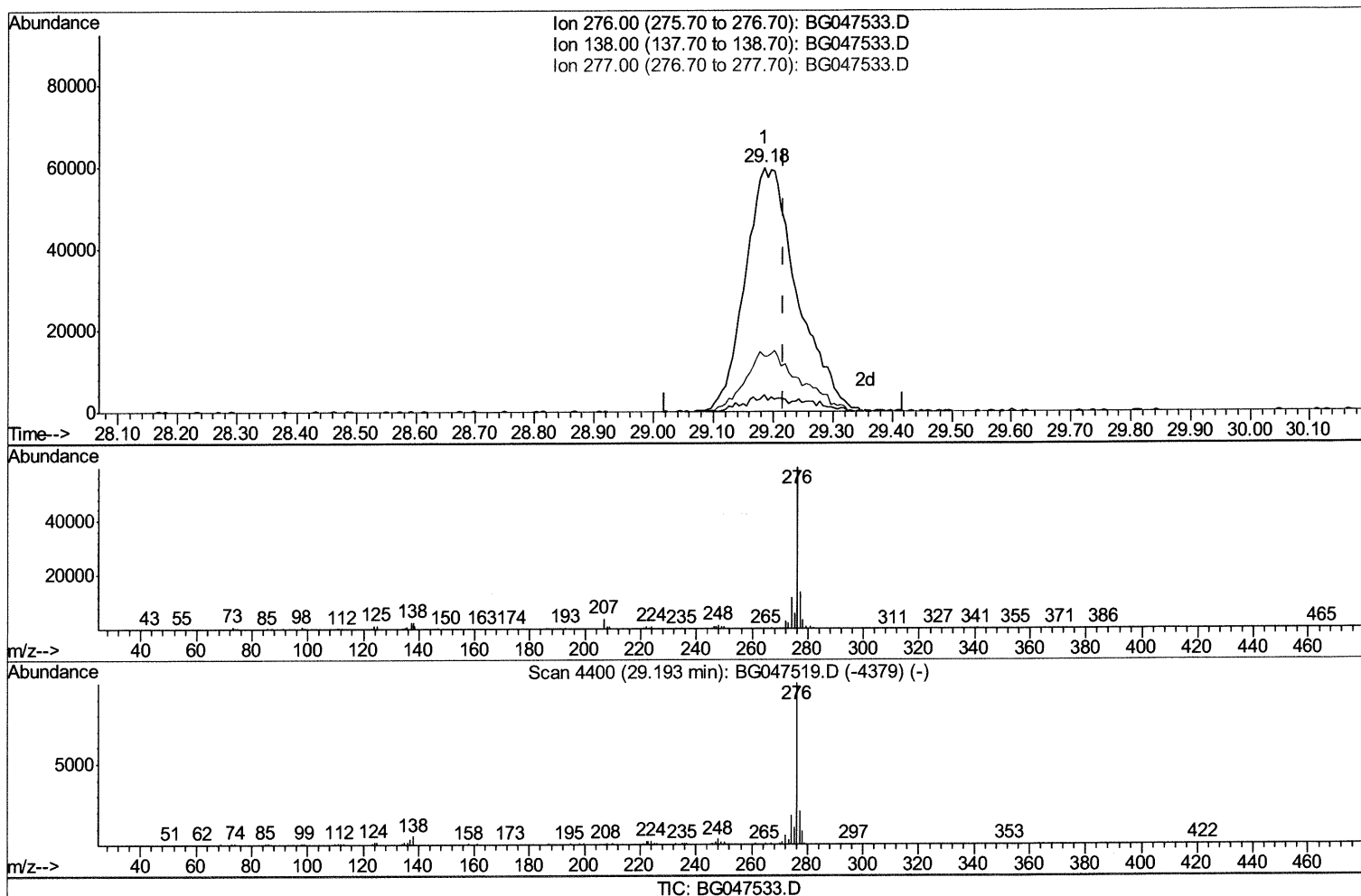
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
Data File : BG047533.D  
Acq On : 2 Dec 2020 16:04  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02069

Manual Integrations  
APPROVED

mohammad  
12/3/2020 3:35:14 PM

Quant Time: Dec 02 16:59:37 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



(92) Indeno(1,2,3-cd)pyrene

29.185min (-0.034) 18.84ng/ul m 12/04/2020

response 360858

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 6.30  | 4.38# |
| 277.00 | 20.60 | 22.82 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

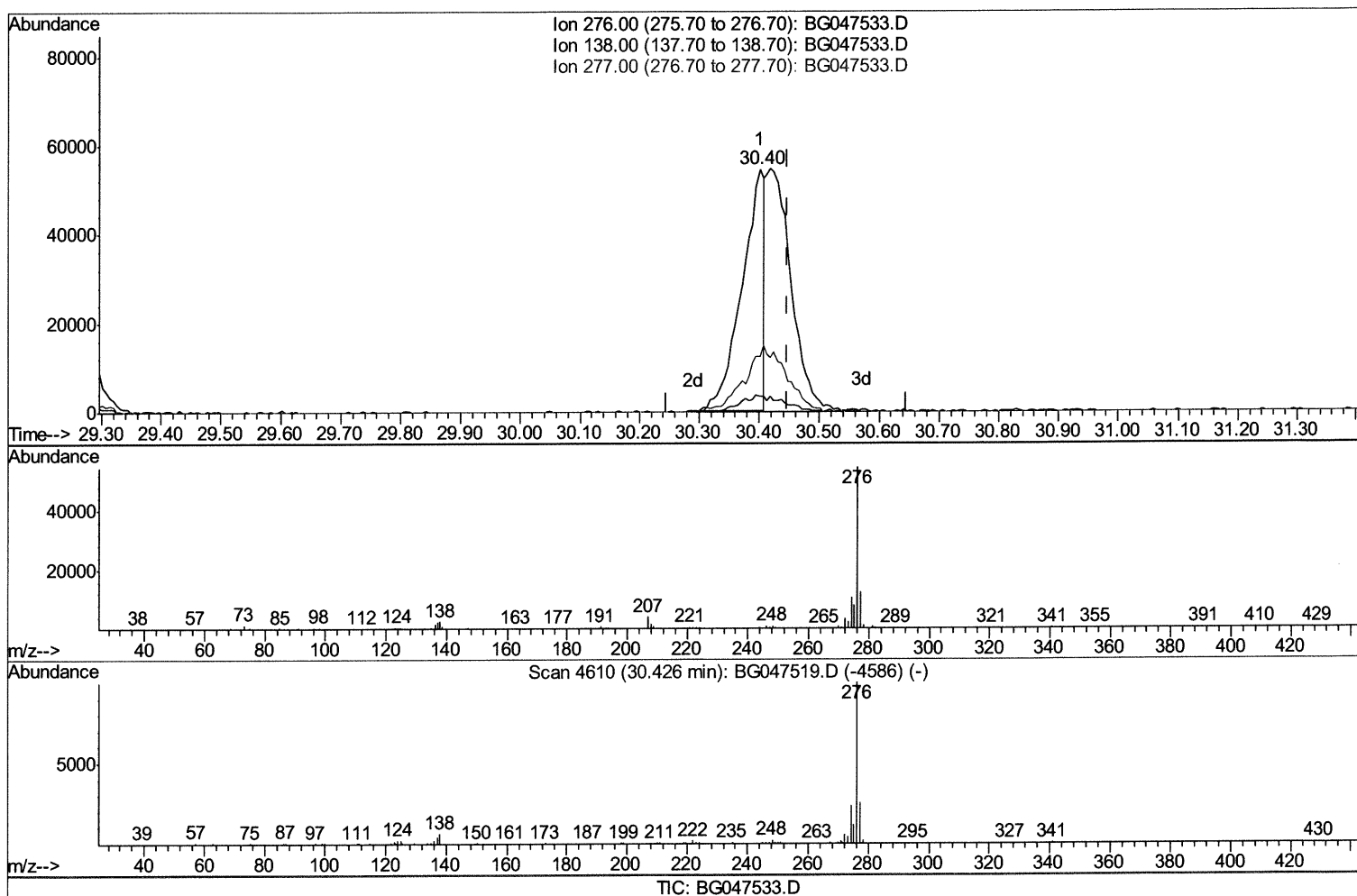
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
Data File : BG047533.D  
Acq On : 2 Dec 2020 16:04  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02069

Manual Integrations  
APPROVED

mohammad  
12/3/2020 3:35:14 PM

Quant Time: Dec 02 16:59:37 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



(94) Benzo(g,h,i)perylene

30.401min (-0.046) 8.77ng/ul

response 138627

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 7.20  | 6.36  |
| 277.00 | 21.80 | 22.97 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

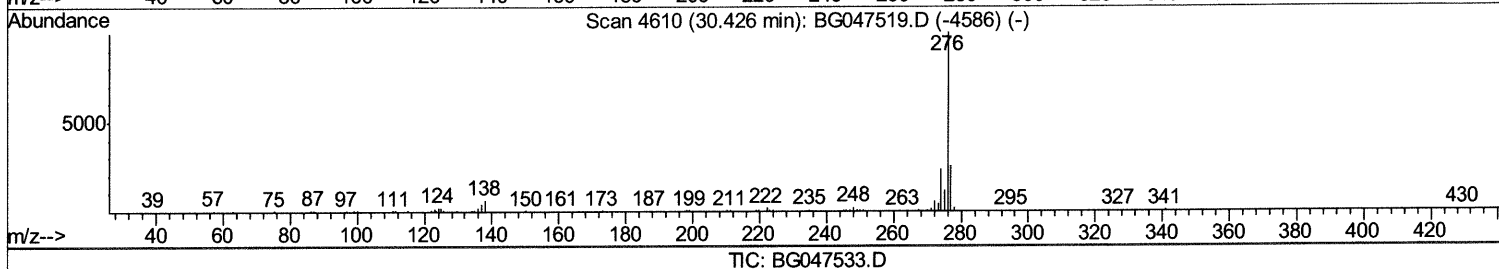
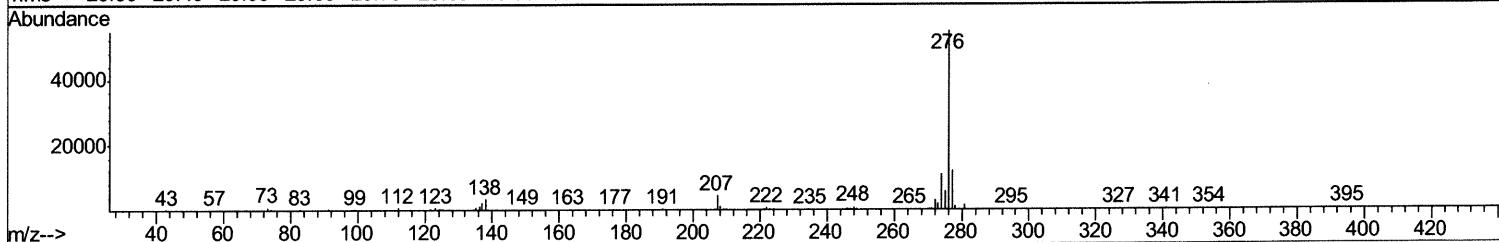
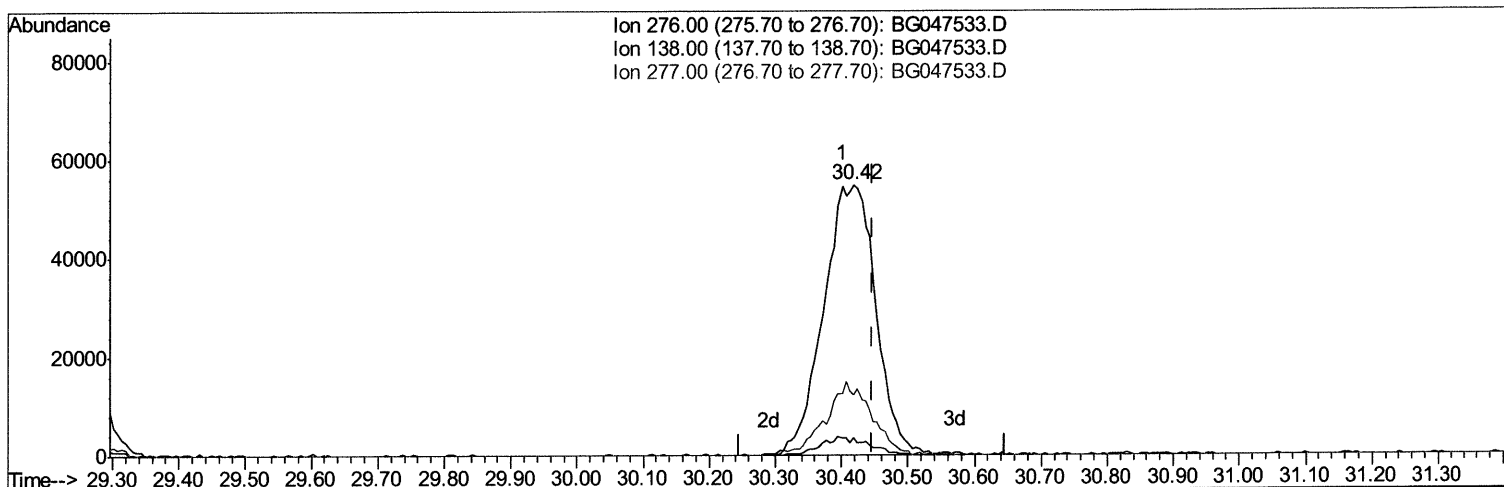
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
Data File : BG047533.D  
Acq On : 2 Dec 2020 16:04  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02069

Manual Integrations  
APPROVED

mohammad  
12/3/2020 3:35:14 PM

Quant Time: Dec 02 16:59:37 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



(94) Benzo(g,h,i)perylene

30.418min (-0.028) 18.79ng/ul m 12/04/20 54

response 297093

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 276.00 | 100   | 100   |
| 138.00 | 7.20  | 6.30  |
| 277.00 | 21.80 | 22.27 |
| 0.00   | 0.00  | 0.00  |

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
 Data File : BG047533.D  
 Acq On : 2 Dec 2020 16:04  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 46 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02069

Manual Integrations  
 APPROVED

mohammad  
 12/3/2020 3:35:14 PM

Quant Time: Dec 02 17:02:46 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

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.27  | 152  | 29652    | 20.00 | ng/ul | -0.01    |
| 18) Naphthalene-d8        | 11.09 | 136  | 119438   | 20.00 | ng/ul | -0.02    |
| 36) Acenaphthene-d10      | 14.88 | 164  | 94690    | 20.00 | ng/ul | -0.01    |
| 62) Phenanthrene-d10      | 17.61 | 188  | 254498   | 20.00 | ng/ul | -0.01    |
| 78) Chrysene-d12          | 21.89 | 240  | 235603   | 20.00 | ng/ul | -0.01    |
| 86) Perylene-d12          | 25.29 | 264  | 250219   | 20.00 | ng/ul | -0.02    |

## System Monitoring Compounds

|                                |       |     |          |       |        |       |             |
|--------------------------------|-------|-----|----------|-------|--------|-------|-------------|
| 3) 1,4-Dioxane-d8              | 3.60  | 96  | 4416m>   | 5.41  | ng/uL> | -0.01 | 12/04/20 JU |
| 5) Phenol-d5                   | 7.40  | 99  | 51379    | 17.28 | ng/ul  | -0.01 |             |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.57  | 67  | 26145    | 15.42 | ng/ul  | -0.01 |             |
| 9) 2-Chlorophenol-d4           | 7.79  | 132 | 36947    | 18.52 | ng/ul  | -0.02 |             |
| 13) 4-Methylphenol-d8          | 8.96  | 113 | 39560    | 17.31 | ng/ul  | -0.01 |             |
| 19) Nitrobenzene-d5            | 9.43  | 128 | 19557    | 19.82 | ng/ul  | -0.01 |             |
| 22) 2-Nitrophenol-d4           | 10.16 | 143 | 23216    | 21.57 | ng/ul  | -0.01 |             |
| 26) 2,4-Dichlorophenol-d3      | 10.69 | 165 | 50412    | 20.45 | ng/ul  | -0.02 |             |
| 29) 4-Chloroaniline-d4         | 11.21 | 131 | 48913    | 19.62 | ng/ul  | -0.02 |             |
| 44) Dimethylphthalate-d6       | 14.27 | 166 | 151492   | 18.64 | ng/ul  | -0.02 |             |
| 47) Acenaphthylene-d8          | 14.57 | 160 | 179315   | 19.09 | ng/ul  | -0.01 |             |
| 52) 4-Nitrophenol-d4           | 15.04 | 143 | 23494    | 18.94 | ng/ul  | 0.00  |             |
| 58) Fluorene-d10               | 15.86 | 176 | 143148   | 19.22 | ng/ul  | -0.01 |             |
| 63) 4,6-Dinitro-2-methylphenol | 15.96 | 200 | 29403    | 20.30 | ng/ul  | -0.01 |             |
| 71) Anthracene-d10             | 17.71 | 188 | 233027m> | 18.46 | ng/ul> | -0.01 | 12/04/20 JU |
| 79) Pyrene-d10                 | 19.98 | 212 | 260432   | 19.00 | ng/ul  | 0.00  |             |
| 90) Benzo(a)pyrene-d12         | 25.05 | 264 | 267068   | 18.54 | ng/ul  | -0.02 |             |

## Target Compounds

|                                |       |     |        |        | Ovalue |    |
|--------------------------------|-------|-----|--------|--------|--------|----|
| 2) 1,4-Dioxane                 | 3.64  | 88  | 3707   | 4.404  | ng/uL# | 62 |
| 4) Benzaldehyde                | 7.39  | 77  | 31783  | 18.095 | ng/ul  | 91 |
| 6) Phenol                      | 7.42  | 94  | 48373  | 16.979 | ng/ul  | 92 |
| 8) Bis(2-Chloroethyl)ether     | 7.67  | 93  | 33844  | 16.093 | ng/ul  | 94 |
| 10) 2-Chlorophenol             | 7.82  | 128 | 36340  | 18.055 | ng/ul# | 88 |
| 11) 2-Methylphenol             | 8.69  | 108 | 37130  | 16.776 | ng/ul  | 95 |
| 12) 2,2'-oxybis(1-Chloropropan | 8.78  | 45  | 33418  | 15.107 | ng/ul# | 93 |
| 14) Acetophenone               | 9.08  | 105 | 64658  | 18.370 | ng/ul  | 93 |
| 15) N-Nitroso-di-n-propylamine | 9.07  | 70  | 35560  | 16.746 | ng/ul# | 87 |
| 16) 4-Methylphenol             | 9.02  | 108 | 40944  | 17.627 | ng/ul  | 99 |
| 17) Hexachloroethane           | 9.36  | 117 | 17988  | 19.018 | ng/ul  | 95 |
| 20) Nitrobenzene               | 9.47  | 77  | 57754  | 18.100 | ng/ul# | 94 |
| 21) Isophorone                 | 10.00 | 82  | 100345 | 16.500 | ng/ul# | 95 |
| 23) 2-Nitrophenol              | 10.19 | 139 | 22698  | 19.396 | ng/ul# | 75 |
| 24) 2,4-Dimethylphenol         | 10.24 | 107 | 54793  | 18.445 | ng/ul  | 94 |
| 25) Bis(2-Chloroethoxy)methane | 10.48 | 93  | 47042  | 16.082 | ng/ul  | 94 |
| 27) 2,4-Dichlorophenol         | 10.72 | 162 | 44825  | 20.485 | ng/ul  | 98 |
| 28) Naphthalene                | 11.14 | 128 | 125326 | 18.526 | ng/ul  | 98 |
| 30) 4-Chloroaniline            | 11.24 | 127 | 48501  | 20.417 | ng/ul  | 99 |
| 31) Hexachlorobutadiene        | 11.42 | 225 | 44517  | 21.624 | ng/ul  | 97 |
| 32) Caprolactam                | 11.99 | 113 | 14861  | 18.633 | ng/ul# | 82 |
| 33) 4-Chloro-3-methylphenol    | 12.33 | 107 | 48646  | 17.988 | ng/ul  | 93 |
| 34) 2-Methylnaphthalene        | 12.73 | 142 | 100370 | 20.022 | ng/ul  | 84 |

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 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 46 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02069

Manual Integrations  
 APPROVED

mohammad  
 12/3/2020 3:35:14 PM

Quant Time: Dec 02 17:02:46 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

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min)    |
|--------------------------------|-------|------|----------|--------|--------|-------------|
| 35) 1-Methylnaphthalene        | 12.94 | 142  | 111666   | 18.952 | ng/uL  | 99          |
| 37) 1,2,4,5-Tetrachlorobenzene | 13.09 | 216  | 74049    | 19.901 | ng/uL  | 98          |
| 38) Hexachlorocyclopentadiene  | 13.06 | 237  | 46250    | 17.863 | ng/uL  | 96          |
| 39) 2,4,6-Trichlorophenol      | 13.31 | 196  | 44731    | 19.359 | ng/uL  | 98          |
| 40) 2,4,5-Trichlorophenol      | 13.39 | 196  | 50577    | 21.083 | ng/uL  | 91          |
| 41) 1,1'-Biphenyl              | 13.71 | 154  | 138552   | 18.864 | ng/uL  | 99          |
| 42) 2-Chloronaphthalene        | 13.76 | 162  | 109696   | 18.841 | ng/uL  | 99          |
| 43) 2-Nitroaniline             | 13.95 | 65   | 40976    | 19.350 | ng/uL  | 95          |
| 45) Dimethylphthalate          | 14.32 | 163  | 156947   | 18.928 | ng/uL  | 97          |
| 46) 2,6-Dinitrotoluene         | 14.44 | 165  | 33082    | 20.835 | ng/uL  | 97          |
| 48) Acenaphthylene             | 14.60 | 152  | 168714   | 19.322 | ng/uL# | 94          |
| 49) 3-Nitroaniline             | 14.77 | 138  | 25186    | 20.313 | ng/uL  | 90          |
| 50) Acenaphthene               | 14.94 | 153  | 115285   | 18.823 | ng/uL  | 98          |
| 51) 2,4-Dinitrophenol          | 14.97 | 184  | 17282    | 20.425 | ng/uL# | 70          |
| 53) 4-Nitrophenol              | 15.05 | 109  | 37715    | 21.004 | ng/uL  | 95          |
| 54) Dibenzofuran               | 15.27 | 168  | 168746   | 18.758 | ng/uL  | 96          |
| 55) 2,4-Dinitrotoluene         | 15.22 | 165  | 47740    | 21.191 | ng/uL  | 93          |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.49 | 232  | 47987    | 20.887 | ng/uL# | 96          |
| 57) Diethylphthalate           | 15.67 | 149  | 153743   | 18.900 | ng/uL  | 99          |
| 59) Fluorene                   | 15.92 | 166  | 144486   | 19.065 | ng/uL  | 94          |
| 60) 4-Chlorophenyl-phenylether | 15.91 | 204  | 91544    | 20.109 | ng/uL  | 93          |
| 61) 4-Nitroaniline             | 15.92 | 138  | 28994    | 19.972 | ng/uL  | 88          |
| 64) 4,6-Dinitro-2-methylphenol | 15.98 | 198  | 30021    | 19.804 | ng/uL# | 97          |
| 65) N-Nitrosodiphenylamine     | 16.11 | 169  | 128876   | 17.741 | ng/uL  | 97          |
| 66) 4-Bromophenyl-phenylether  | 16.79 | 248  | 62714    | 19.398 | ng/uL  | 93          |
| 67) Hexachlorobenzene          | 16.92 | 284  | 68367    | 20.215 | ng/uL  | 96          |
| 68) Atrazine                   | 17.05 | 200  | 59374    | 18.569 | ng/uL  | 94          |
| 69) Pentachlorophenol          | 17.26 | 266  | 31628    | 15.988 | ng/uL  | 90          |
| 70) Phenanthrene               | 17.66 | 178  | 249933   | 18.186 | ng/uL  | 99          |
| 72) Anthracene                 | 17.75 | 178  | 252982   | 18.278 | ng/uL  | 98          |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.69 | 216  | 76303    | 19.073 | ng/uL  | 96          |
| 74) Pentachlorobenzene         | 15.19 | 250  | 73323    | 19.063 | ng/uL  | 95          |
| 75) Carbazole                  | 18.00 | 167  | 201566   | 18.054 | ng/uL  | 96          |
| 76) Di-n-butylphthalate        | 18.55 | 149  | 251805   | 17.706 | ng/uL  | 98          |
| 77) Fluoranthene               | 19.64 | 202  | 329987   | 19.012 | ng/uL  | 99          |
| 80) Perylene                   | 20.00 | 202  | 317866   | 19.046 | ng/uL  | 99          |
| 81) Butylbenzylphthalate       | 20.87 | 149  | 105662   | 18.027 | ng/uL  | 95          |
| 82) 3,3'-Dichlorobenzidine     | 21.78 | 252  | 102639   | 18.661 | ng/uL  | 85          |
| 83) Benzo(a)anthracene         | 21.88 | 228  | 303949   | 18.132 | ng/uL  | 99          |
| 84) Bis(2-ethylhexyl)phthalate | 21.76 | 149  | 147216   | 18.019 | ng/uL  | 95          |
| 85) Chrysene                   | 21.94 | 228  | 294599   | 18.467 | ng/uL  | 99          |
| 87) Di-n-octyl phthalate       | 23.03 | 149  | 244845   | 17.668 | ng/uL  | 100         |
| 88) Benzo(b)fluoranthene       | 24.21 | 252  | 336790   | 19.063 | ng/uL  | 97          |
| 89) Benzo(k)fluoranthene       | 24.28 | 252  | 314045   | 18.038 | ng/uL  | 98          |
| 91) Benzo(a)pyrene             | 25.13 | 252  | 292836   | 18.594 | ng/uL# | 99          |
| 92) Indeno(1,2,3-cd)pyrene     | 29.18 | 276  | 360858m> | 18.837 | ng/uL> | 12/04/20 J4 |
| 93) Dibenzo(a,h)anthracene     | 29.26 | 278  | 304194   | 19.367 | ng/uL# | 94          |
| 94) Benzo(a,h,i)perylene       | 30.42 | 276  | 297093m> | 18.795 | ng/uL> | 12/04/20 J4 |

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120120\  
Data File : BG047533.D  
Acq On : 2 Dec 2020 16:04  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02069

Manual Integrations  
APPROVED

mohammad  
12/3/2020 3:35:14 PM

Quant Time: Dec 02 17:02:46 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

| Internal Standards                                                         | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                      |      |      |          |      |       |          |
| (#) = qualifier out of range (m) = manual integration (+) = signals summed |      |      |          |      |       |          |