

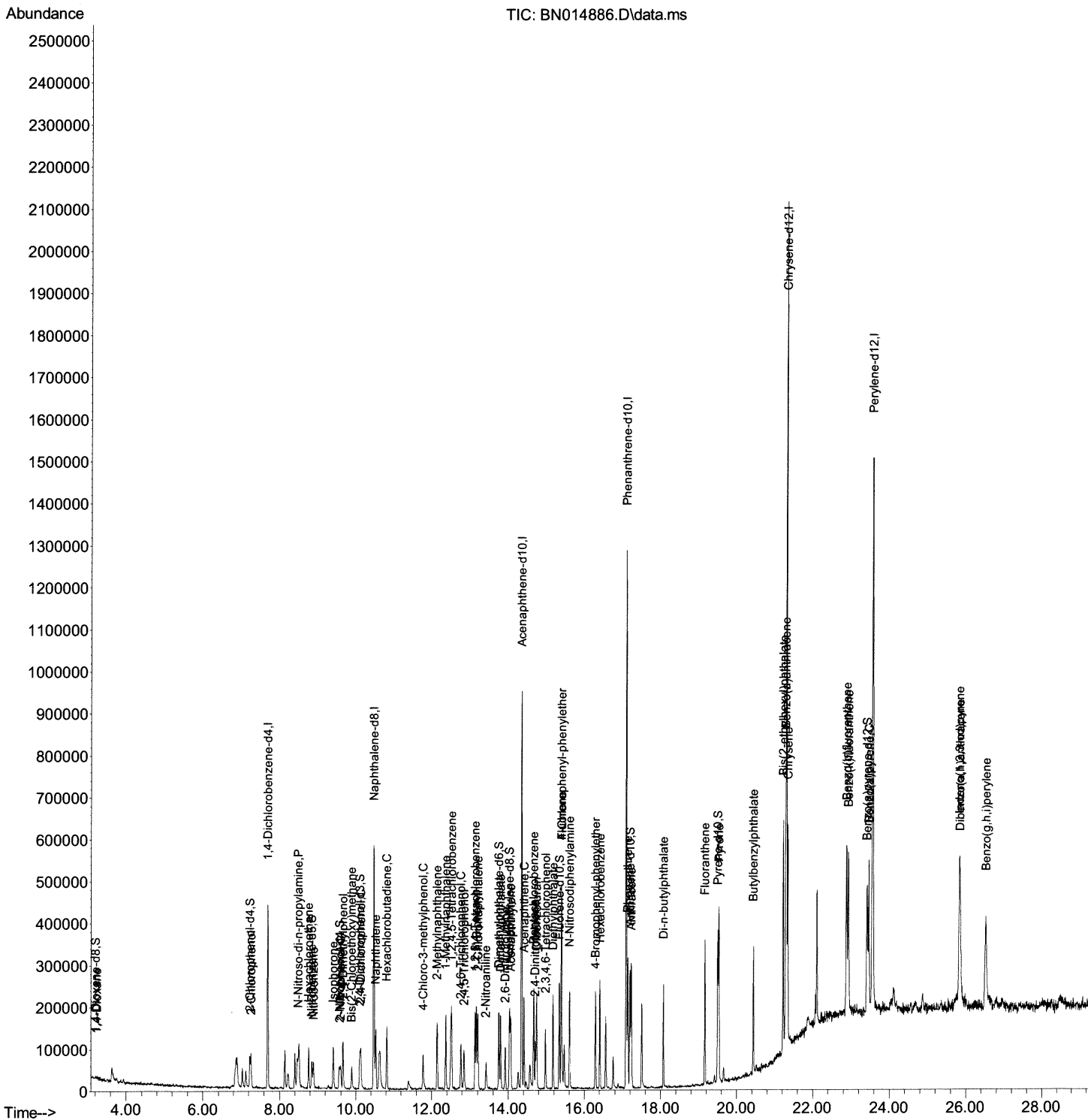
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
Data File : BN014886.D  
Acq On : 12 Jun 2021 10:02  
Operator : CG/JU  
Sample : SST00551  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD005051

Manual Integrations  
APPROVED

mohammad  
6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jun 14 01:48:30 2021  
Response via : Initial Calibration



# Quantitation Report (Qedit)

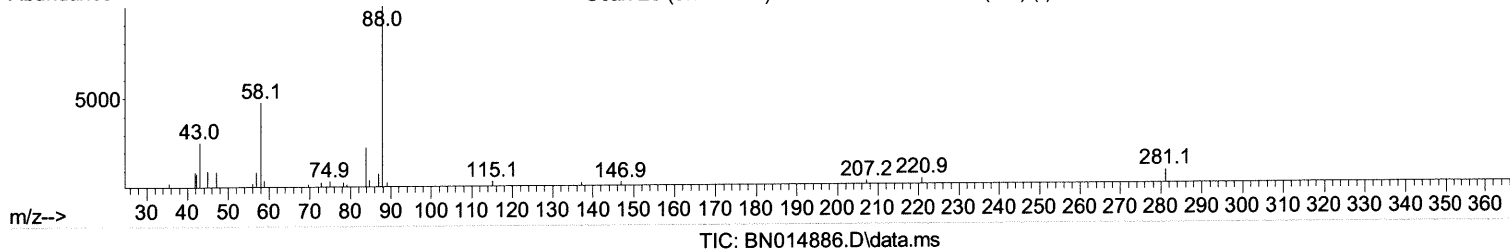
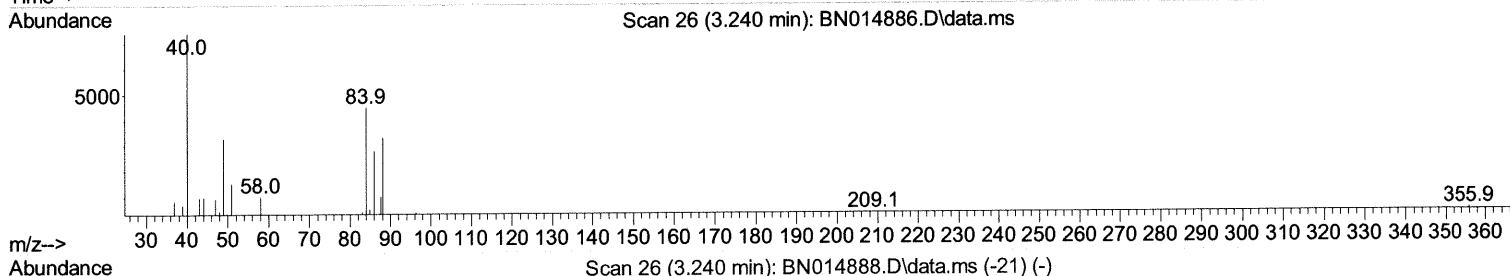
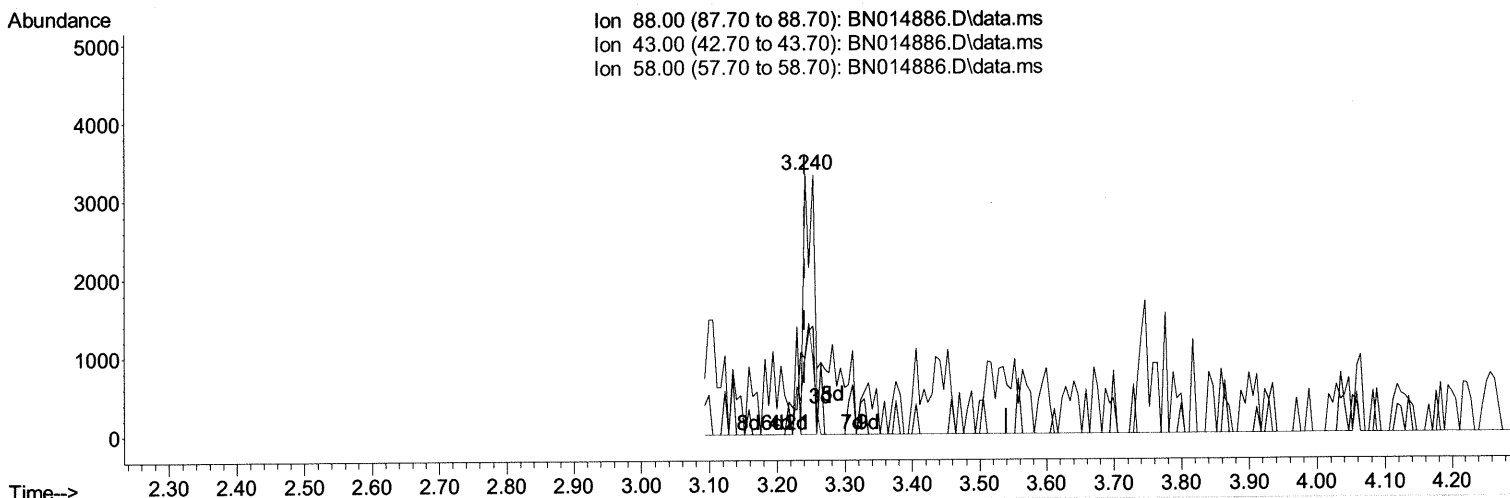
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SSTD00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD005051

Manual Integrations  
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mohammad  
 6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



(2) 1,4-Dioxane

3.240min (+ 0.000) 1.17 ng/uL

response 3384

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 88.00 | 100.00 | 100.00 |
| 43.00 | 17.20  | 29.08# |
| 58.00 | 46.40  | 30.49# |
| 0.00  | 0.00   | 0.00   |

## Quantitation Report (Qedit)

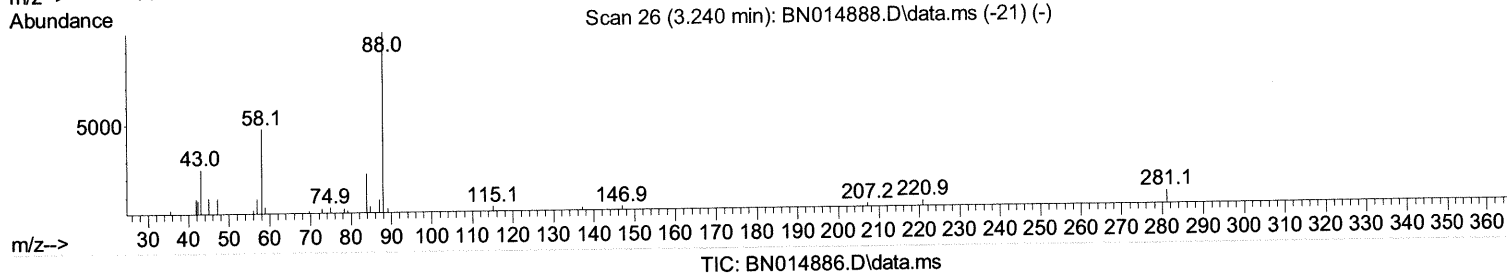
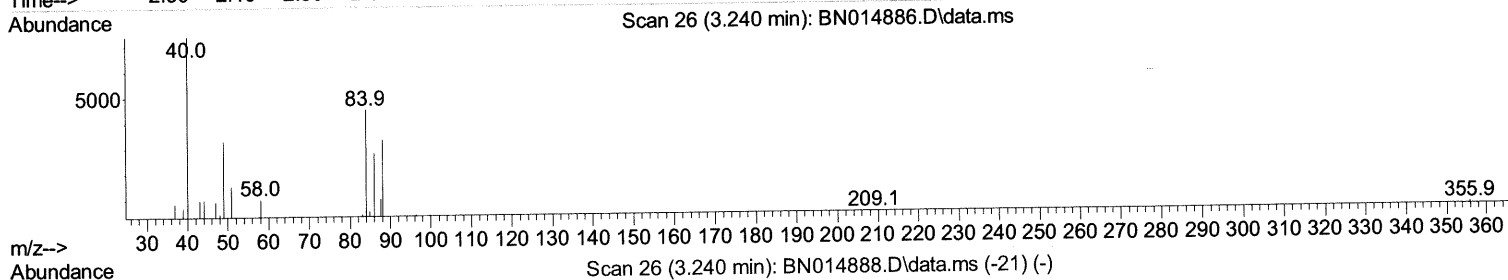
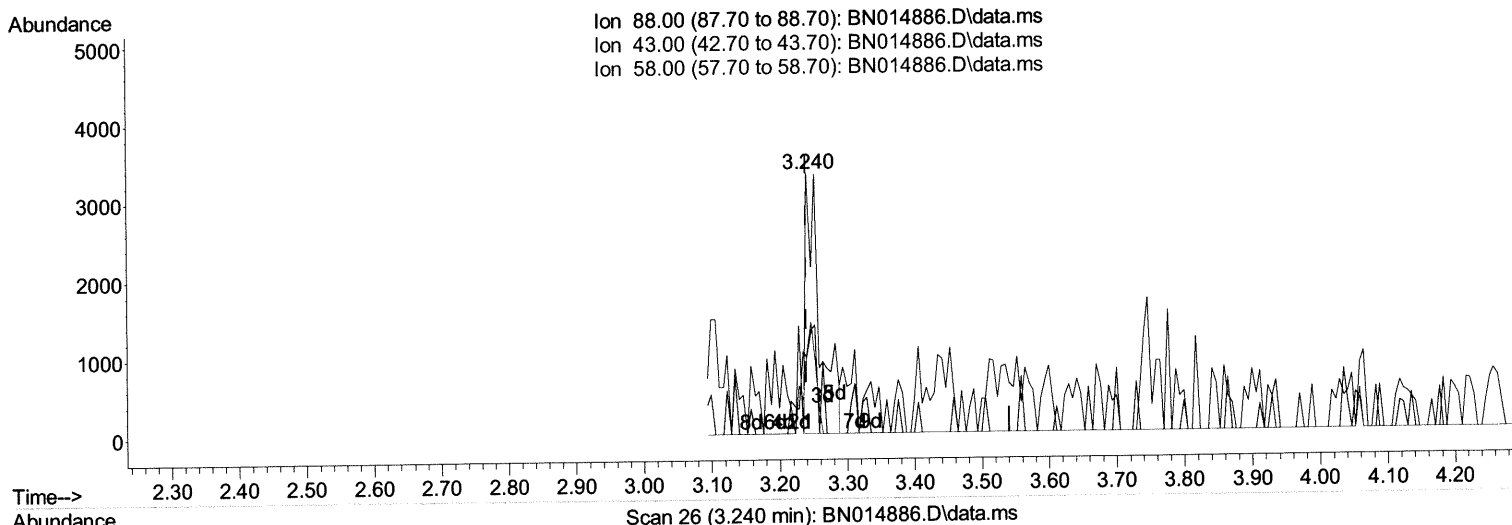
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Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061221\  
Data File : BN014886.D  
Acq On    : 12 Jun 2021  10:02  
Operator  : CG/JU  
Sample    : SSTD00551  
Misc      :  
ALS Vial  : 3    Sample Multiplier: 1
```

**Instrument :**  
BNA\_N  
**ClientSampleId :**  
SSTD005051

**Manual Integrations**  
**APPROVED**

mohammad  
6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jun 14 01:48:30 2021  
Response via : Initial Calibration



(2) 1,4-Dioxane

3.240min (+ 0.000) 1.69 ng/uL m 06/15/21 JY

response 4896

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 88.00 | 100.00 | 100.00 |
| 43.00 | 17.20  | 29.08# |
| 58.00 | 46.40  | 30.49# |
| 0.00  | 0.00   | 0.00   |

# Quantitation Report (Qedit)

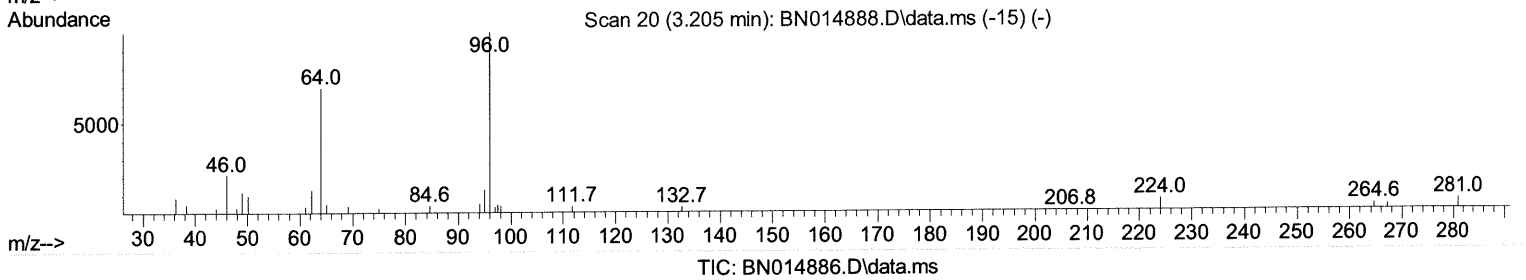
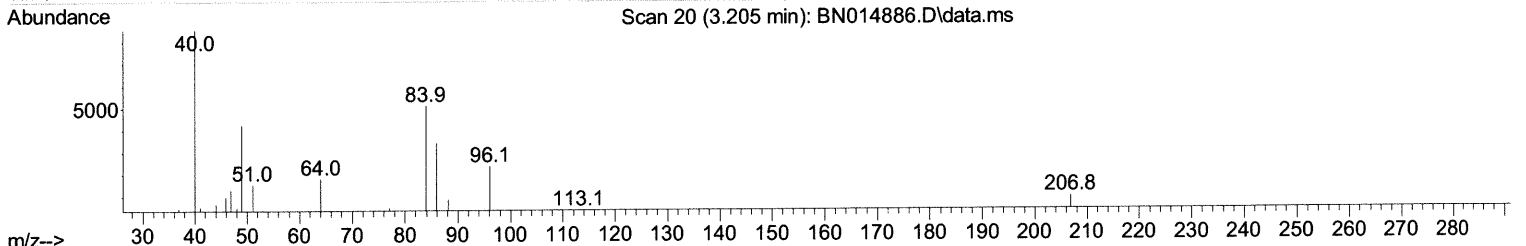
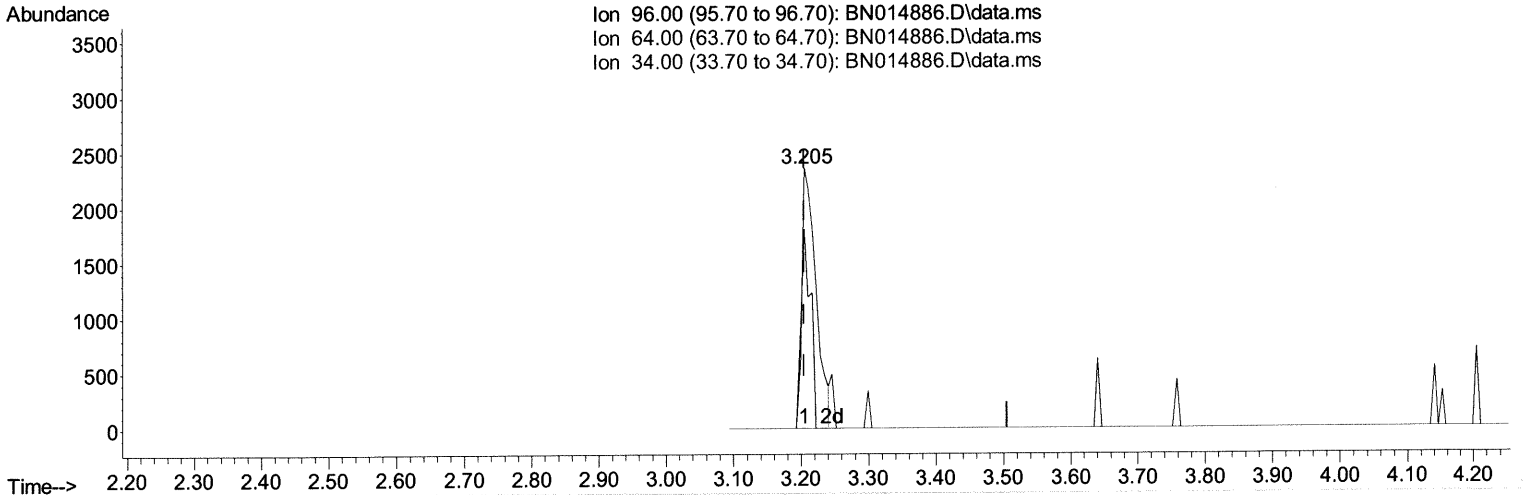
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SSTD00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD005051

Manual Integrations  
 APPROVED

mohammad  
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Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



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

3.205min (+ 0.000) 1.40 ng/uL

response 3445

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 64.30  | 76.62  |
| 34.00 | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

# Quantitation Report (Qedit)

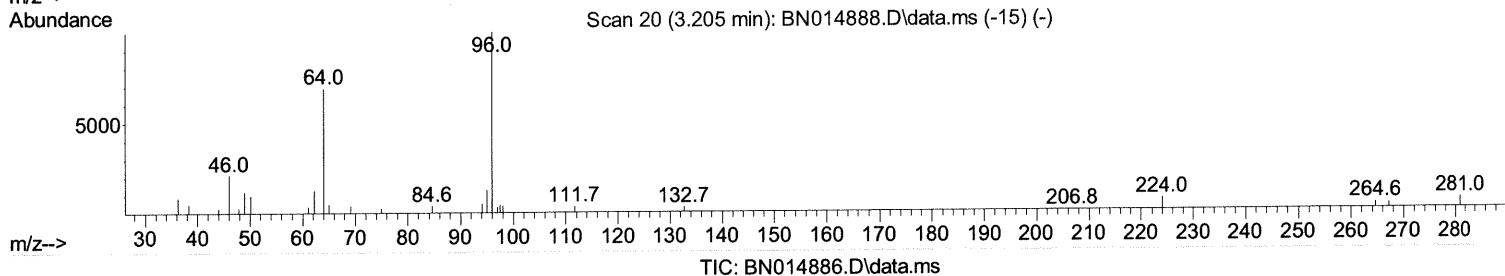
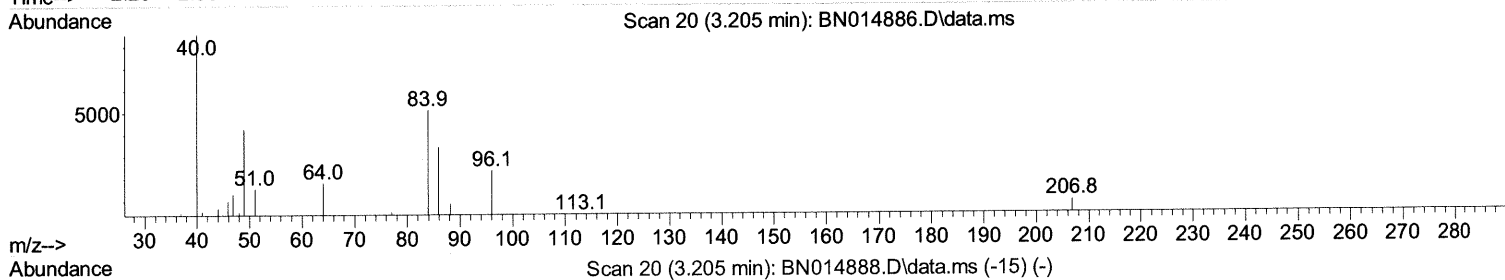
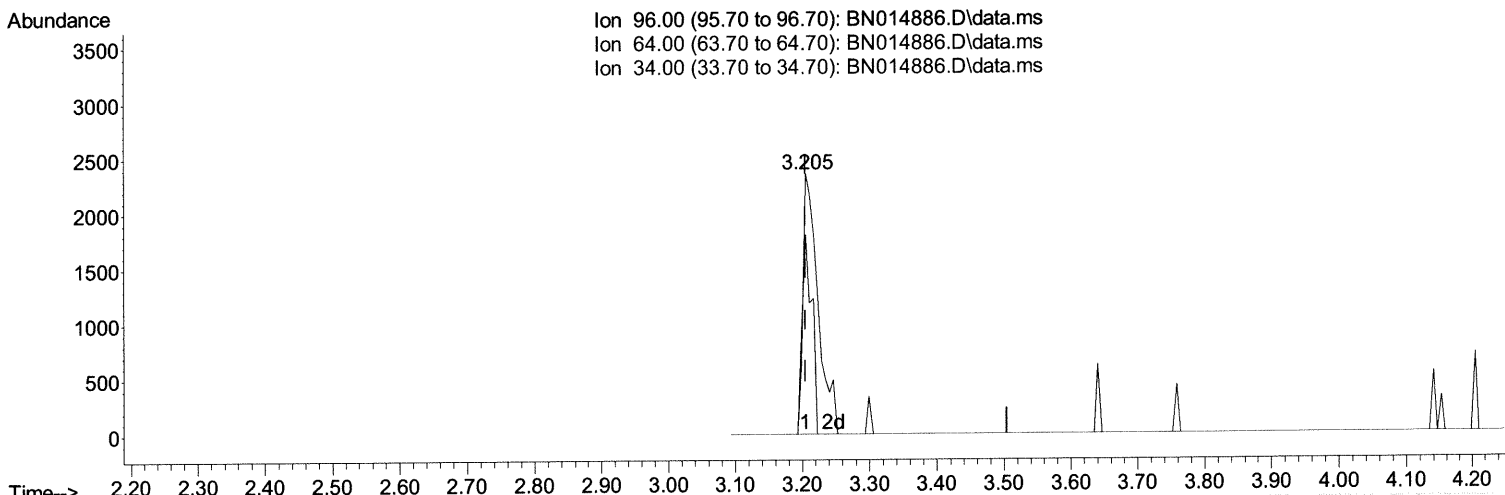
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SSTD00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD005051

Manual Integrations  
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Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



TIC: BN014886.D\data.ms

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

3.205min (+ 0.000) 1.47 ng/uL m 06/15/21 JU

response 3616

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 64.30  | 76.62  |
| 34.00 | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

# Quantitation Report (Qedit)

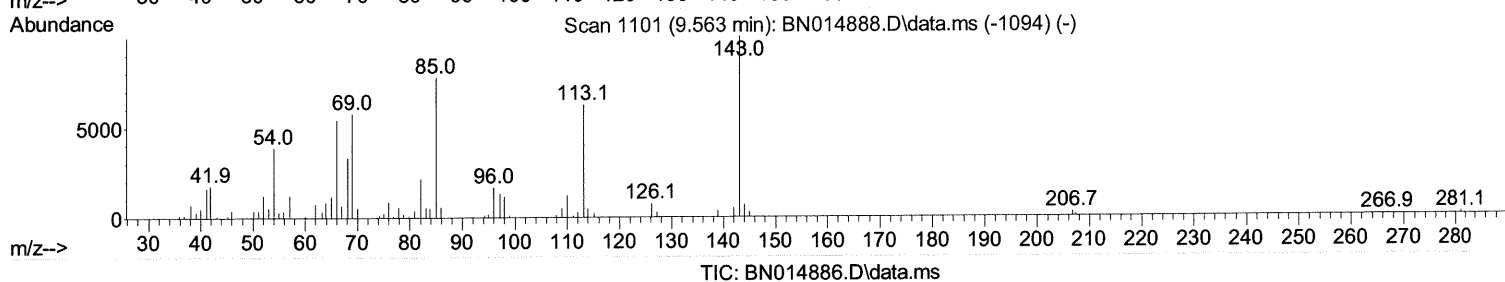
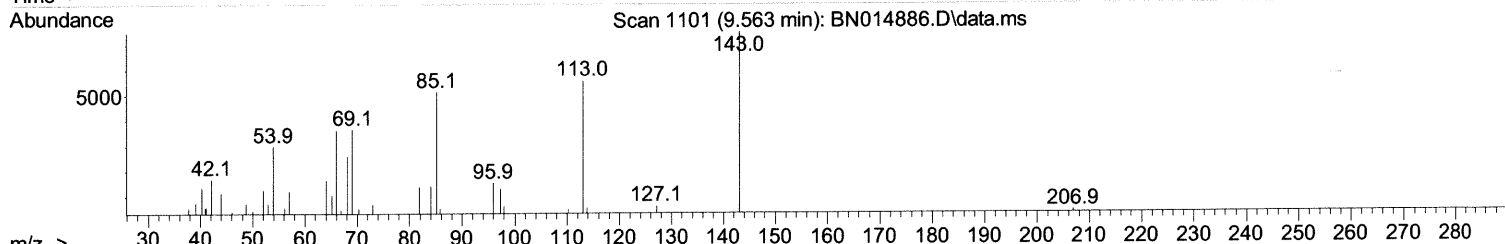
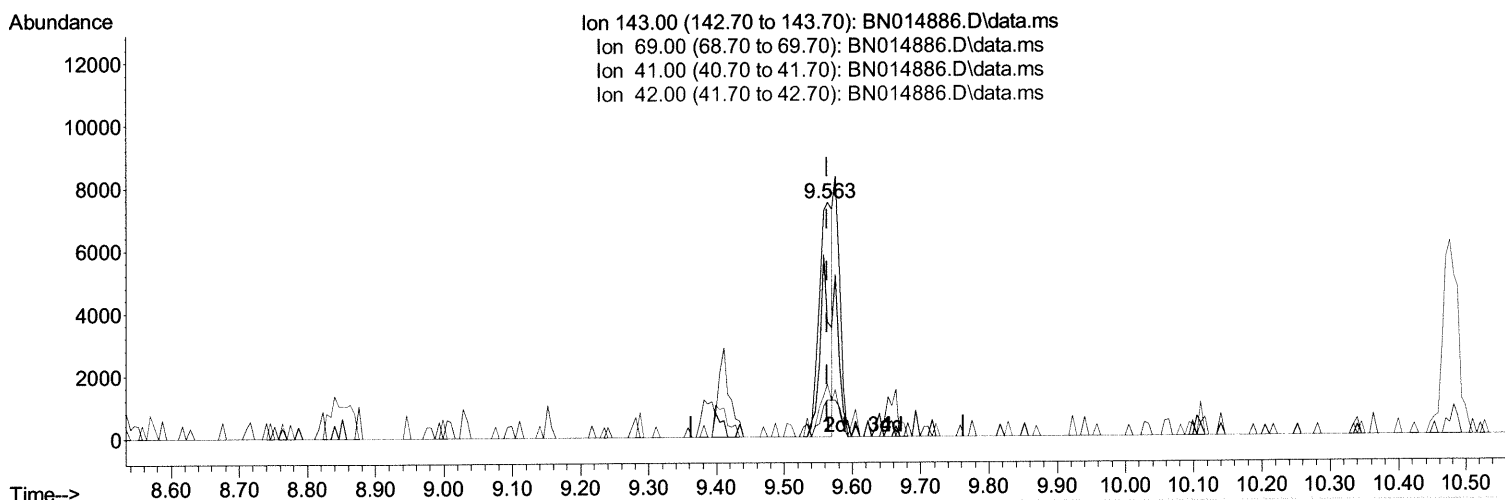
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SSTD00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD005051

Manual Integrations  
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mohammad  
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Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



(24) 2-Nitrophenol-d4 (S)

9.563min (+ 0.000) 3.14 ng/ul

response 10113

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 143.00 | 100.00 | 100.00 |
| 69.00  | 69.60  | 48.98# |
| 41.00  | 18.30  | 15.54  |
| 42.00  | 14.10  | 22.52# |

## Quantitation Report (Qedit)

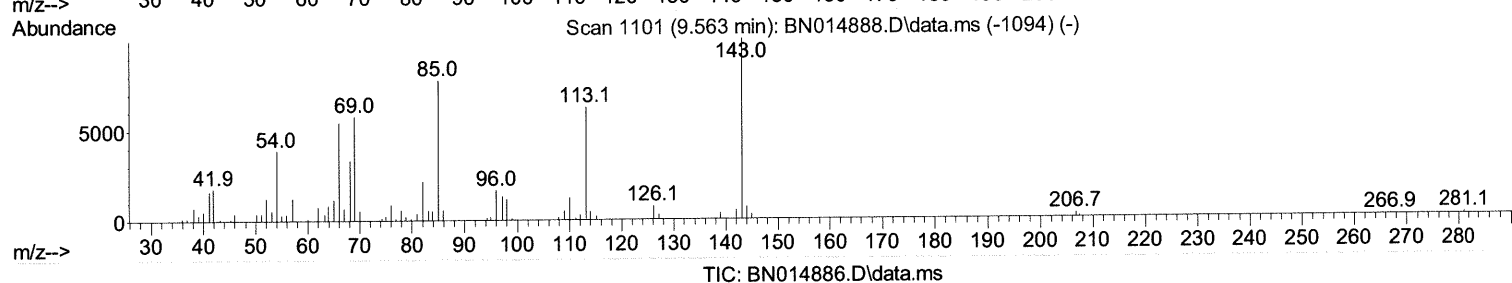
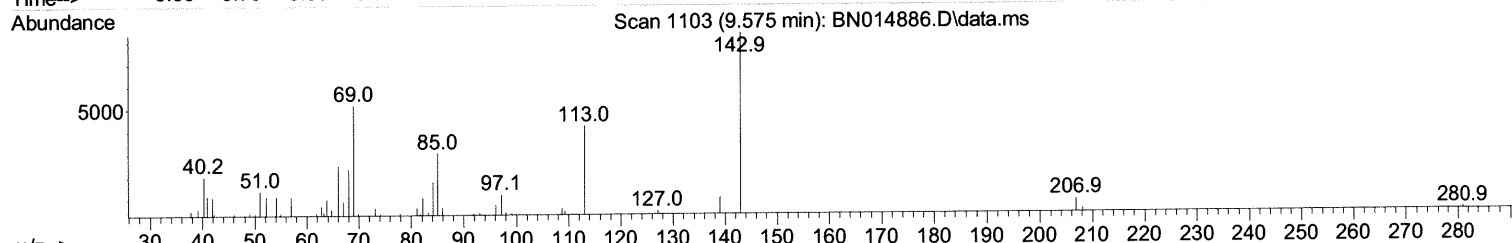
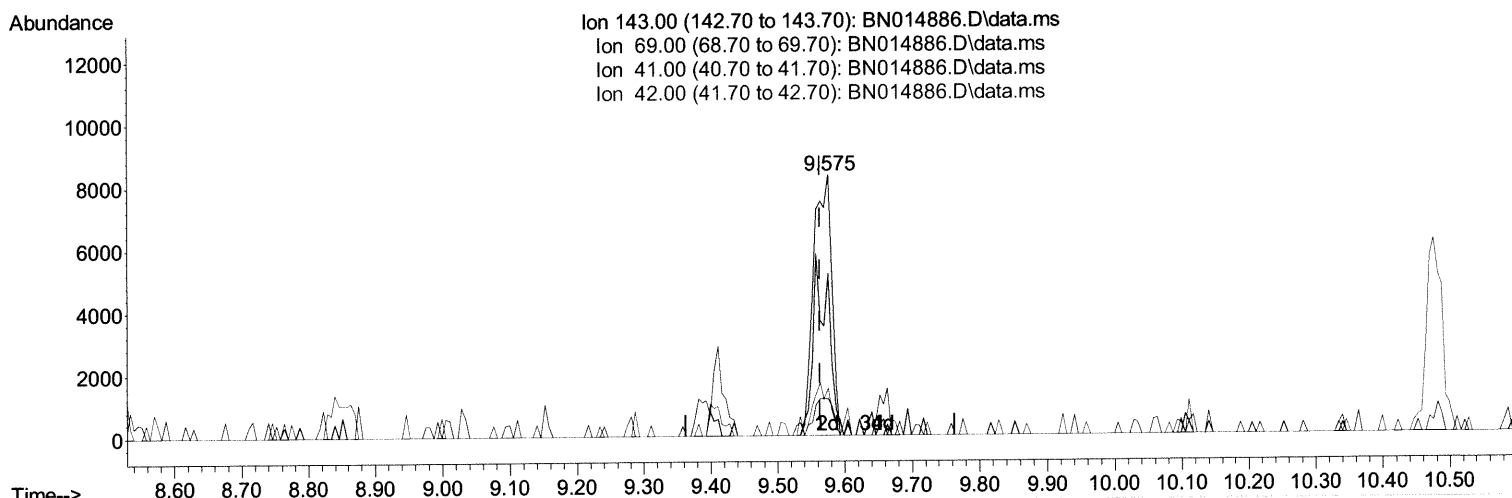
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
Data File : BN014886.D  
Acq On : 12 Jun 2021 10:02  
Operator : CG/JU  
Sample : SSTD00551  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD005051

Manual Integrations  
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Quant Time: Jun 14 01:50:03 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jun 14 01:48:30 2021  
Response via : Initial Calibration



TIC: BN014886.D\data.ms

(24) 2-Nitrophenol-d4 (S)

9.575min (+ 0.012) 4.78 ng/ul m 06/15/21 JU

response 15402

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 143.00 | 100.00 | 100.00 |
| 69.00  | 69.60  | 62.16  |
| 41.00  | 18.30  | 14.01# |
| 42.00  | 14.10  | 13.36  |

# Quantitation Report (Qedit)

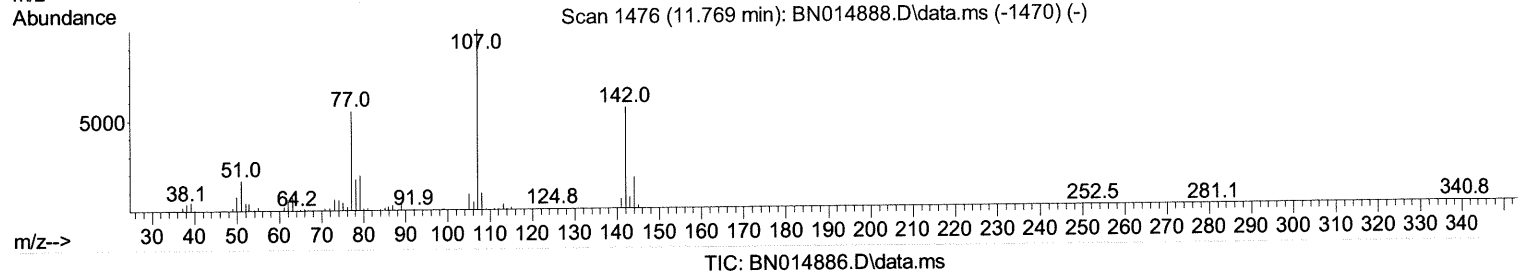
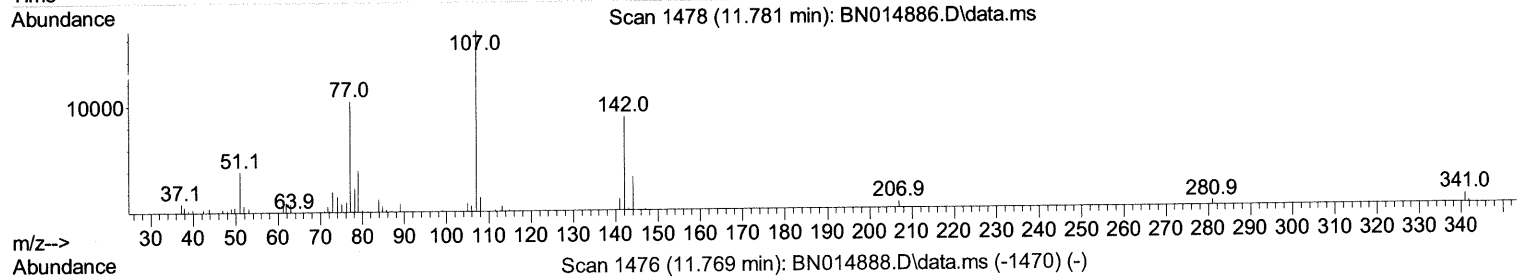
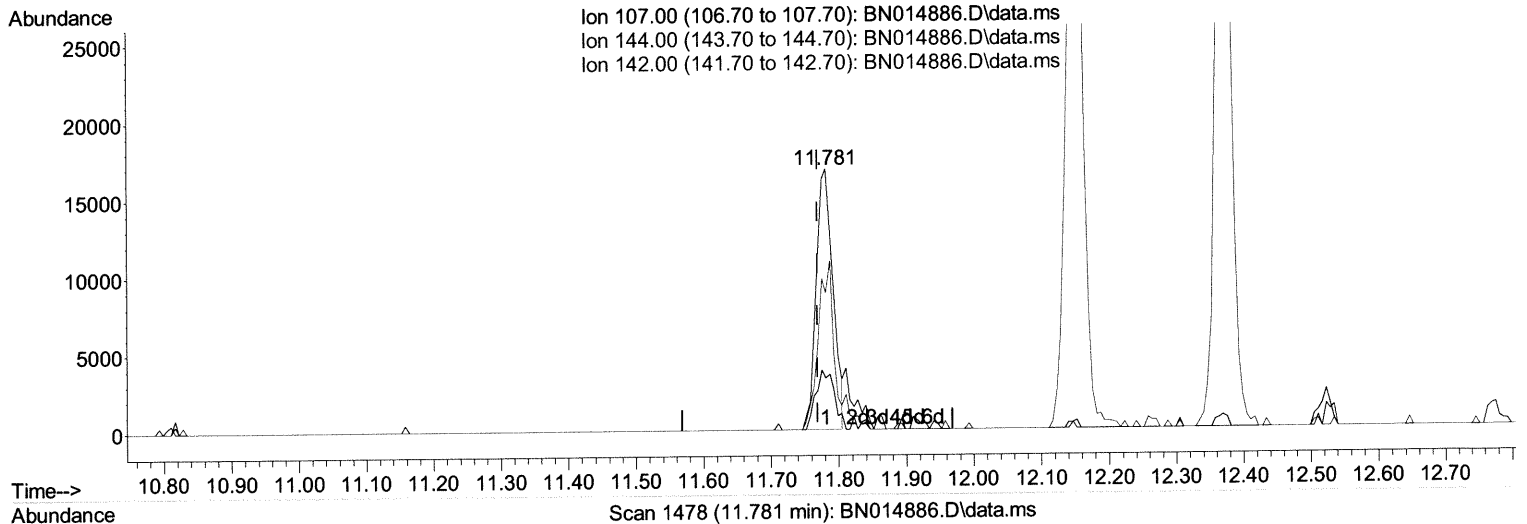
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SSTD00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD005051

Manual Integrations  
 APPROVED

mohammad  
 6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



TIC: BN014886.D\data.ms

(35) 4-Chloro-3-methylphenol (C)

11.781min (+ 0.012) 2.85 ng/ul

response 29085

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 107.00 | 100.00 | 100.00 |
| 144.00 | 16.80  | 19.68  |
| 142.00 | 53.20  | 52.54  |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

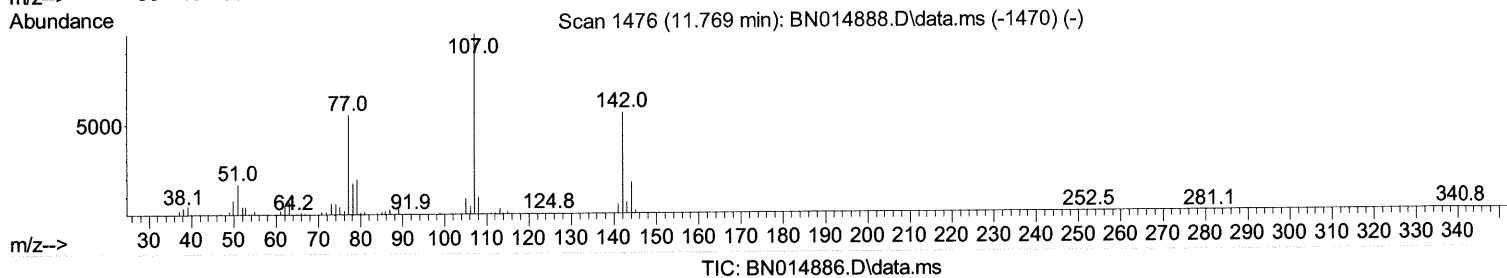
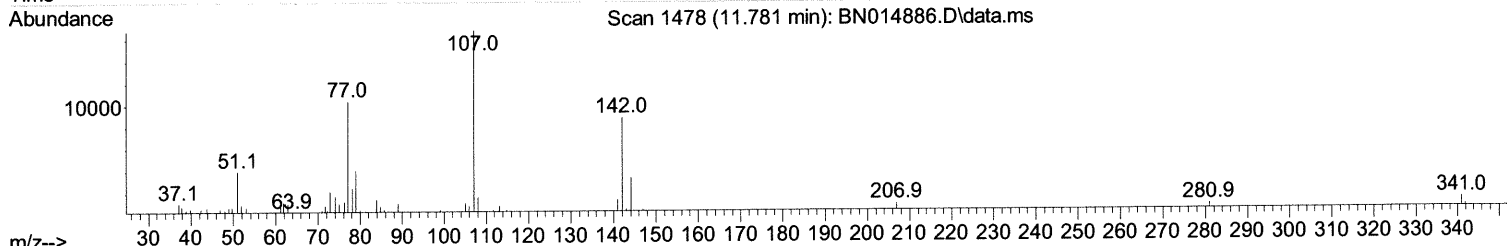
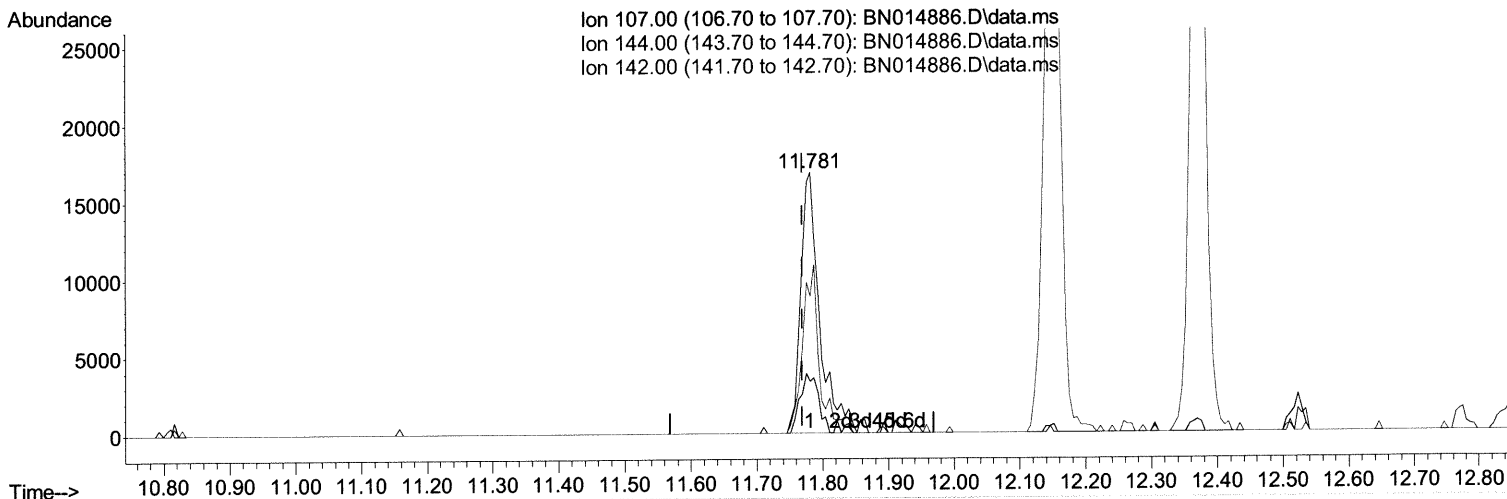
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SST00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST005051

Manual Integrations  
 APPROVED

mohammad  
 6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



(35) 4-Chloro-3-methylphenol (C)

11.781min (+ 0.012) 3.27 ng/ul m 06/15/21 JU

response 33419

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 107.00 | 100.00 | 100.00 |
| 144.00 | 16.80  | 19.68  |
| 142.00 | 53.20  | 52.54  |
| 0.00   | 0.00   | 0.00   |

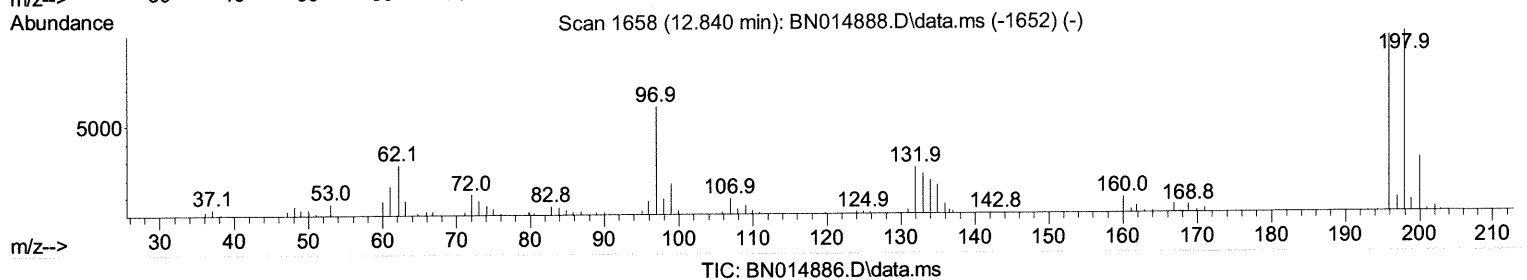
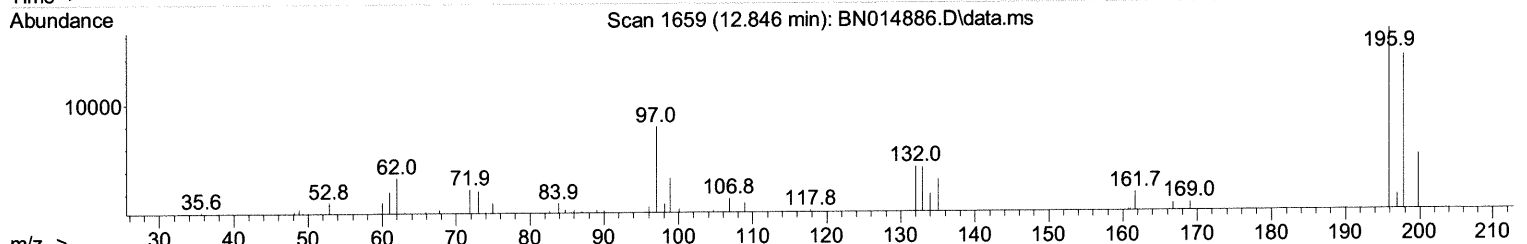
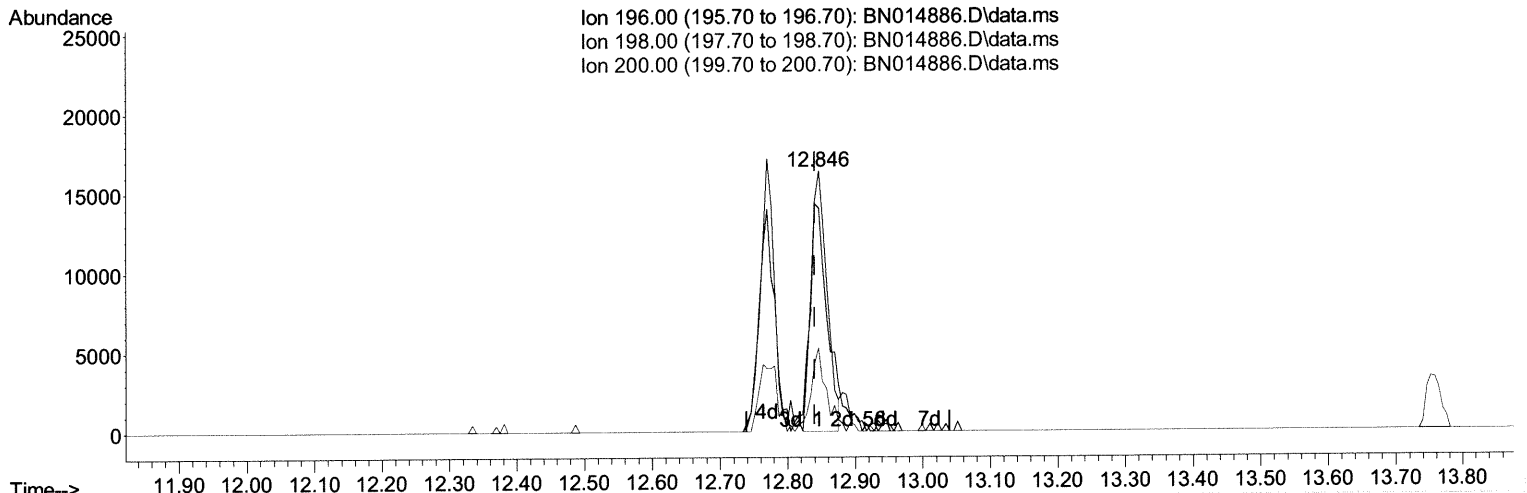
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Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061221\  
Data File : BN014886.D  
Acq On    : 12 Jun 2021  10:02  
Operator  : CG/JU  
Sample    : SSTD00551  
Misc      :  
ALS Vial  : 3    Sample Multiplier: 1
```

**Instrument :**  
BNA\_N  
**ClientSampleId :**  
SSTD005051

## Manual Integrations APPROVED

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6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jun 14 01:48:30 2021  
Response via : Initial Calibration



TIC: BN014886.D\data.ms

(42) 2,4,5-Trichlorophenol

12.846min (+ 0.006) 3.96 ng/ul

**response**      26890

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 196.00 | 100.00 | 100.00 |
| 198.00 | 97.30  | 85.62  |
| 200.00 | 30.40  | 31.90  |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

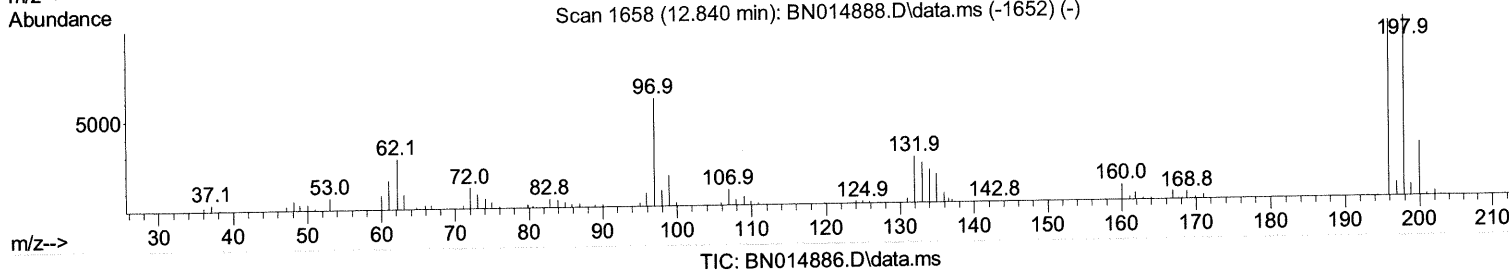
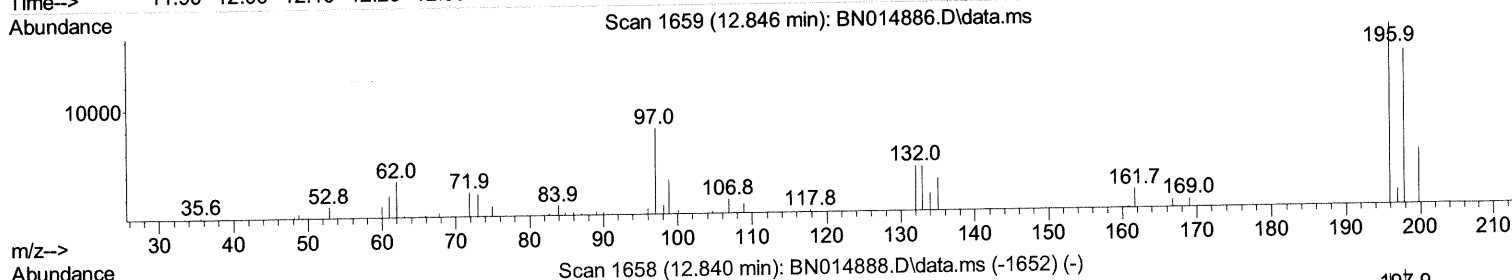
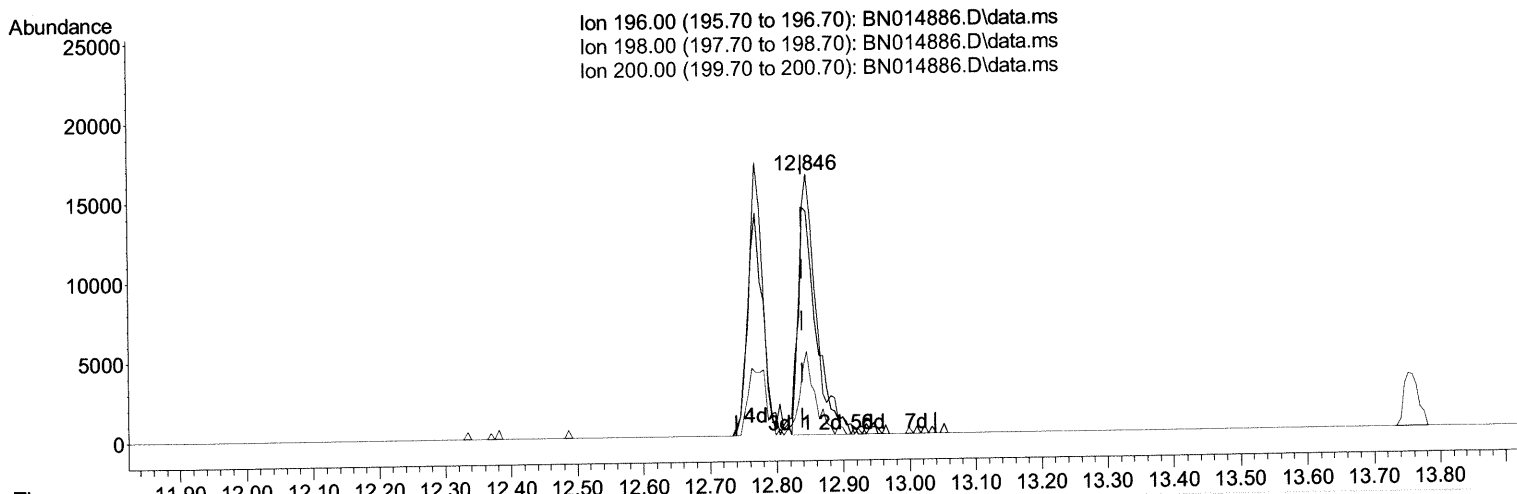
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SSTD00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD005051

Manual Integrations  
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mohammad  
 6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



(42) 2,4,5-Trichlorophenol

12.846min (+ 0.006) 4.37 ng/ul m 06/15/21 JU

response 29662

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 196.00 | 100.00 | 100.00 |
| 198.00 | 97.30  | 85.62  |
| 200.00 | 30.40  | 31.90  |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

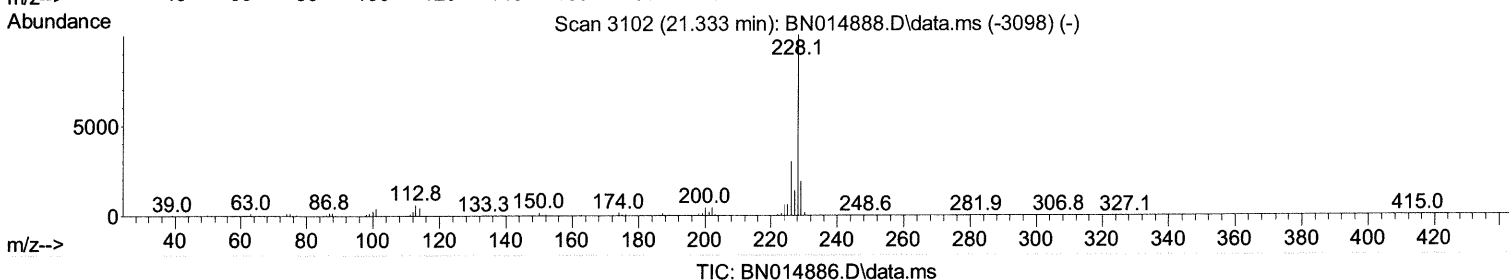
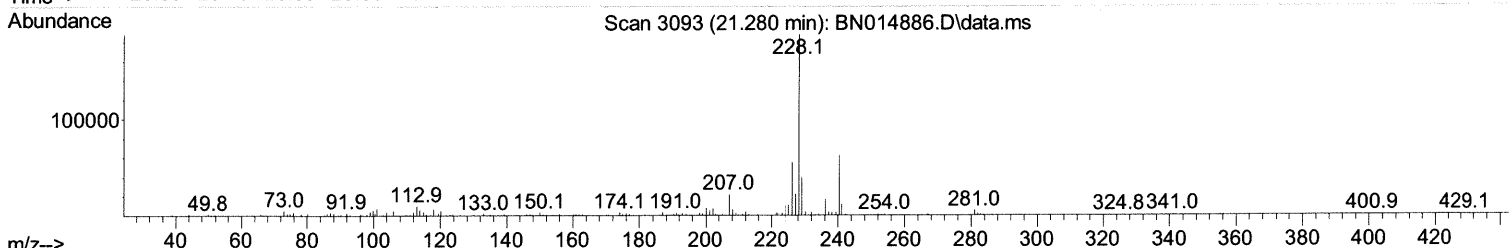
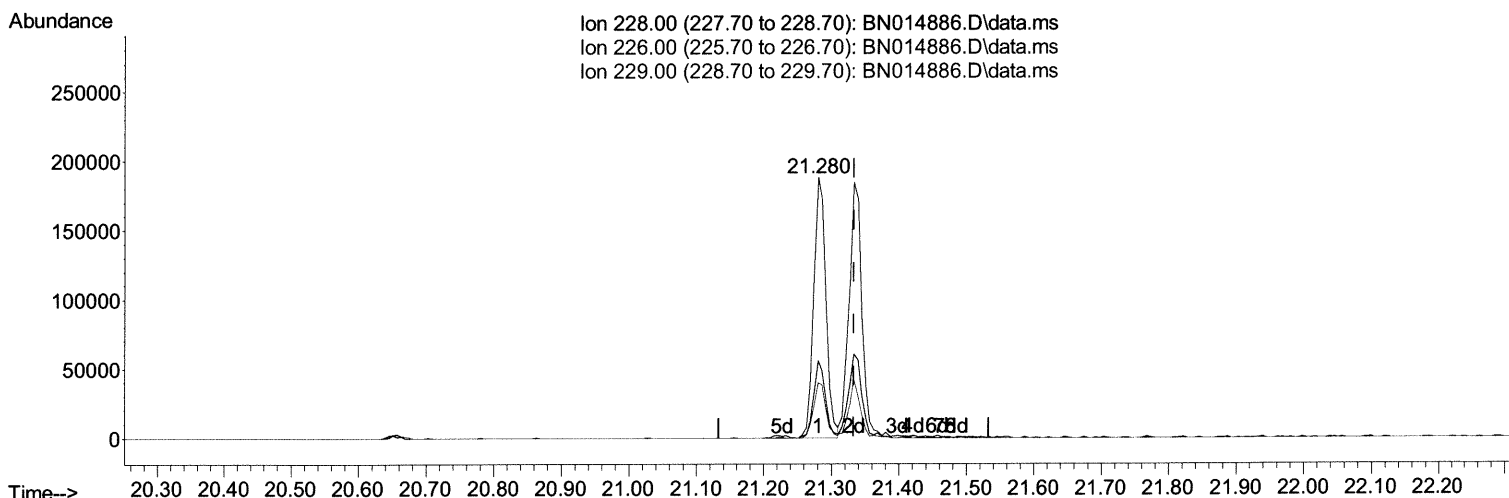
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SST00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST005051

Manual Integrations  
 APPROVED

mohammad  
 6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



(87) Chrysene

21.280min (-0.053) 4.26 ng/ul

response 237251

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 228.00 | 100.00 | 100.00 |
| 226.00 | 29.70  | 29.70  |
| 229.00 | 18.50  | 21.06  |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

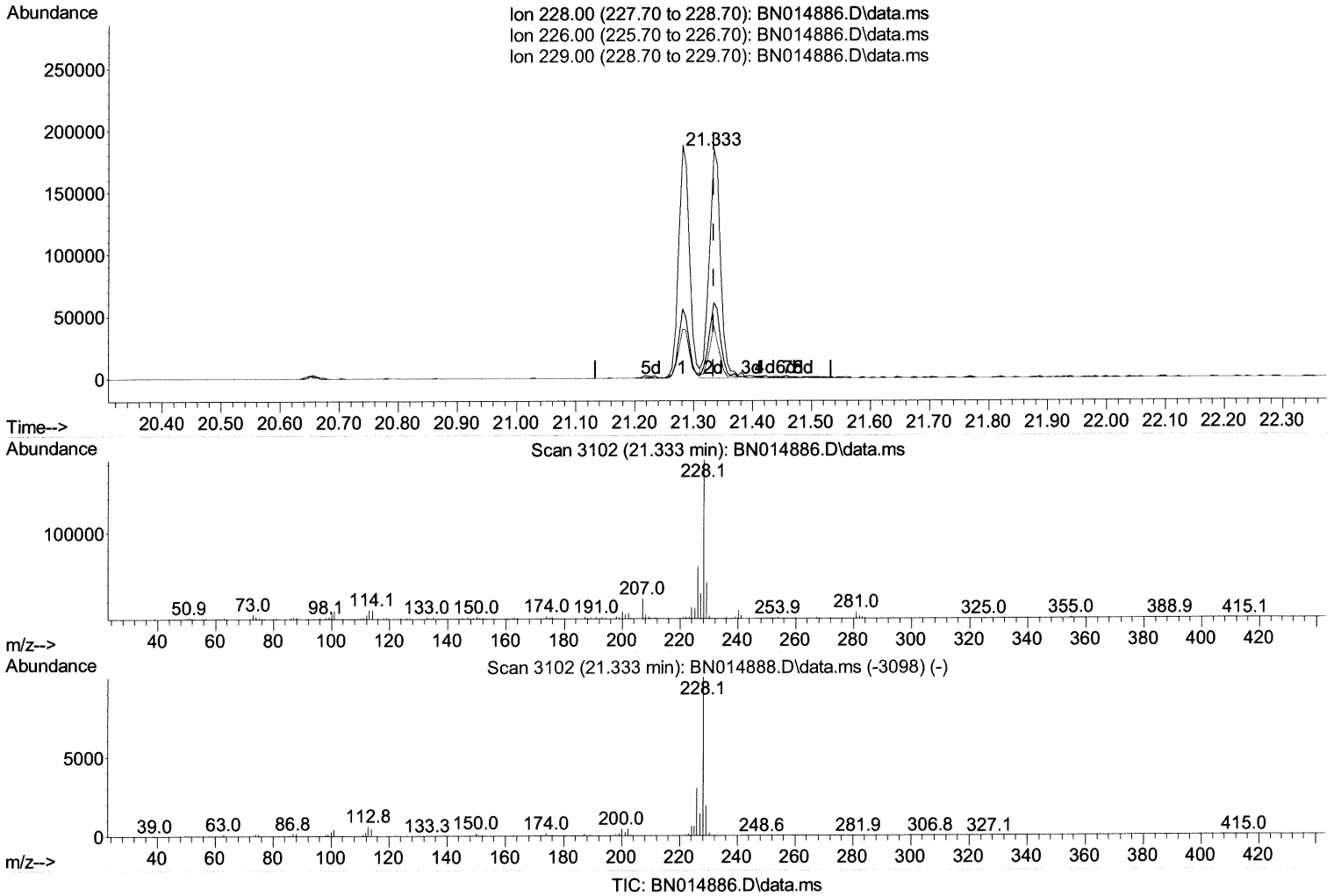
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SSTD00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD005051

Manual Integrations  
 APPROVED

mohammad  
 6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration



(87) Chrysene

21.333min (+ 0.000) 4.38 ng/ul m 06/15/21 JU

response 243413

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 228.00 | 100.00 | 100.00 |
| 226.00 | 29.70  | 32.85  |
| 229.00 | 18.50  | 22.98# |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SST00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampled :  
 SST005051

Manual Integrations  
 APPROVED

mohammad  
 6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min)           |
|-------------------------------|--------|------|----------|--------|--------|--------------------|
| Internal Standards            |        |      |          |        |        |                    |
| 1) 1,4-Dichlorobenzene-d4     | 7.693  | 152  | 100304   | 20.000 | ng/ul  | 0.00               |
| 20) Naphthalene-d8            | 10.475 | 136  | 398171   | 20.000 | ng/ul  | # 0.00             |
| 38) Acenaphthene-d10          | 14.346 | 164  | 293229   | 20.000 | ng/ul  | 0.00               |
| 64) Phenanthrene-d10          | 17.098 | 188  | 696852   | 20.000 | ng/ul  | # 0.00             |
| 79) Chrysene-d12              | 21.298 | 240  | 886255   | 20.000 | ng/ul  | 0.00               |
| 88) Perylene-d12              | 23.551 | 264  | 969481   | 20.000 | ng/ul  | 0.00               |
| System Monitoring Compounds   |        |      |          |        |        |                    |
| 3) 1,4-Dioxane-d8             | 3.205  | 96   | 3616m    | 1.468  | ng/ul  | > 0.00 06/15/21 JU |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/ul  |                    |
| 7) Phenol-d5                  | 0.000  | 99   | 0d       | 0.000  | ng/ul  |                    |
| 9) Bis-(2-Chloroethyl)eth...  | 0.000  | 67   | 0d       | 0.000  | ng/ul  |                    |
| 11) 2-Chlorophenol-d4         | 7.228  | 132  | 22853    | 3.757  | ng/ul  | 0.00               |
| 15) 4-Methylphenol-d8         | 0.000  | 113  | 0d       | 0.000  | ng/ul  |                    |
| 21) Nitrobenzene-d5           | 8.846  | 128  | 11842    | 3.948  | ng/ul  | 0.00               |
| 24) 2-Nitrophenol-d4          | 9.575  | 143  | 15402m   | 4.777  | ng/ul  | > 0.01 06/15/21 JU |
| 28) 2,4-Dichlorophenol-d3     | 10.110 | 165  | 27341    | 3.719  | ng/ul  | 0.01               |
| 31) 4-Chloroaniline-d4        | 0.000  | 131  | 0d       | 0.000  | ng/ul  |                    |
| 46) Dimethylphthalate-d6      | 13.757 | 166  | 110117   | 4.727  | ng/ul  | 0.00               |
| 49) Acenaphthylene-d8         | 14.034 | 160  | 118879   | 4.413  | ng/ul  | 0.00               |
| 54) 4-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000  | ng/ul  |                    |
| 60) Fluorene-d10              | 15.340 | 176  | 90619    | 4.514  | ng/ul  | 0.00               |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200  | 0d       | 0.000  | ng/ul  |                    |
| 73) Anthracene-d10            | 17.198 | 188  | 145459   | 4.468  | ng/ul  | 0.00               |
| 81) Pyrene-d10                | 19.498 | 212  | 187313   | 4.482  | ng/ul  | 0.00               |
| 92) Benzo(a)pyrene-d12        | 23.410 | 264  | 227692   | 4.409  | ng/ul  | 0.00               |
| Target Compounds              |        |      |          |        |        |                    |
| 2) 1,4-Dioxane                | 3.240  | 88   | 4896m    | 1.693  | ng/ul  | > 0.00 06/15/21 JU |
| 12) 2-Chlorophenol            | 7.264  | 128  | 28864    | 4.534  | ng/ul# | 68                 |
| 17) N-Nitroso-di-n-propyla... | 8.493  | 70   | 21759    | 3.574  | ng/ul# | 83                 |
| 19) Hexachloroethane          | 8.763  | 117  | 15822    | 4.964  | ng/ul# | 81                 |
| 22) Nitrobenzene              | 8.893  | 77   | 34439    | 3.679  | ng/ul# | 97                 |
| 23) Isophorone                | 9.405  | 82   | 64736    | 3.642  | ng/ul# | 97                 |
| 25) 2-Nitrophenol             | 9.599  | 139  | 14315    | 4.178  | ng/ul# | 85                 |
| 26) 2,4-Dimethylphenol        | 9.658  | 107  | 41003    | 3.888  | ng/ul# | 74                 |
| 27) Bis(2-Chloroethoxy)met... | 9.893  | 93   | 37415    | 3.464  | ng/ul# | 96                 |
| 29) 2,4-Dichlorophenol        | 10.134 | 162  | 26225    | 3.720  | ng/ul# | 87                 |
| 30) Naphthalene               | 10.528 | 128  | 95885    | 4.431  | ng/ul  | 98                 |
| 33) Hexachlorobutadiene       | 10.816 | 225  | 30115    | 4.777  | ng/ul  | 98                 |
| 35) 4-Chloro-3-methylphenol   | 11.781 | 107  | 33419m   | 3.271  | ng/ul  | > 0.00 06/15/21 JU |
| 36) 2-Methylnaphthalene       | 12.146 | 142  | 69697    | 4.320  | ng/ul  | 92                 |
| 37) 1-Methylnaphthalene       | 12.369 | 142  | 71526    | 4.510  | ng/ul# | 99                 |
| 39) 1,2,4,5-Tetrachloroben... | 12.522 | 216  | 46815    | 4.223  | ng/ul# | 86                 |
| 41) 2,4,6-Trichlorophenol     | 12.769 | 196  | 24589    | 3.950  | ng/ul# | 80                 |
| 42) 2,4,5-Trichlorophenol     | 12.846 | 196  | 29662m   | 4.366  | ng/ul  | > 0.00 06/15/21 JU |
| 43) 1,1'-Biphenyl             | 13.175 | 154  | 97337    | 4.640  | ng/ul# | 97                 |
| 44) 2-Chloronaphthalene       | 13.216 | 162  | 76932    | 4.577  | ng/ul  | 88                 |
| 45) 2-Nitroaniline            | 13.422 | 65   | 18470    | 3.608  | ng/ul  | 89                 |
| 47) Dimethylphthalate         | 13.804 | 163  | 100864   | 4.377  | ng/ul  | 97                 |
| 48) 2,6-Dinitrotoluene        | 13.922 | 165  | 18214    | 4.254  | ng/ul# | 84                 |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061221\  
 Data File : BN014886.D  
 Acq On : 12 Jun 2021 10:02  
 Operator : CG/JU  
 Sample : SSTD00551  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD005051

Manual Integrations  
 APPROVED

mohammad  
 6/14/2021 5:49:15 PM

Quant Time: Jun 14 01:50:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN061221MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 14 01:48:30 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| 50) Acenaphthylene            | 14.063 | 152  | 108947   | 4.399 | ng/ul  | 96       |
| 52) Acenaphthene              | 14.410 | 153  | 77420    | 4.314 | ng/ul  | 89       |
| 56) Dibenzofuran              | 14.746 | 168  | 119998   | 4.495 | ng/ul  | 94       |
| 57) 2,4-Dinitrotoluene        | 14.710 | 165  | 27648    | 4.052 | ng/ul# | 94       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.981 | 232  | 25359    | 3.628 | ng/ul# | 95       |
| 59) Diethylphthalate          | 15.175 | 149  | 103883   | 4.690 | ng/ul  | 97       |
| 61) Fluorene                  | 15.398 | 166  | 103482   | 5.152 | ng/ul  | 93       |
| 62) 4-Chlorophenyl-phenyle... | 15.398 | 204  | 53314    | 4.487 | ng/ul# | 73       |
| 67) N-Nitrosodiphenylamine    | 15.610 | 169  | 82506    | 4.396 | ng/ul  | 98       |
| 68) 4-Bromophenyl-phenylether | 16.298 | 248  | 36100    | 3.978 | ng/ul# | 78       |
| 69) Hexachlorobenzene         | 16.410 | 284  | 56665    | 4.992 | ng/ul# | 91       |
| 72) Phenanthrene              | 17.140 | 178  | 166369   | 4.569 | ng/ul# | 93       |
| 74) Anthracene                | 17.234 | 178  | 163179   | 4.447 | ng/ul# | 95       |
| 75) 1,2,3,4-Tetrachloroben... | 13.140 | 216  | 49852    | 4.329 | ng/ul# | 86       |
| 76) Pentachlorobenzene        | 14.669 | 250  | 57739    | 4.848 | ng/ul# | 91       |
| 78) Di-n-butylphthalate       | 18.075 | 149  | 160343   | 4.400 | ng/ul# | 96       |
| 80) Fluoranthene              | 19.163 | 202  | 207808   | 4.217 | ng/ul# | 94       |
| 82) Pyrene                    | 19.528 | 202  | 239158   | 4.720 | ng/ul# | 92       |
| 83) Butylbenzylphthalate      | 20.439 | 149  | 77460    | 4.549 | ng/ul# | 78       |
| 85) Benzo(a)anthracene        | 21.280 | 228  | 237251   | 4.295 | ng/ul  | 96       |
| 86) Bis(2-ethylhexyl)phtha... | 21.222 | 149  | 116342   | 4.740 | ng/ul# | 89       |
| 87) Chrysene                  | 21.333 | 228  | 243413m> | 4.375 | ng/ul> | 94       |
| 90) Benzo(b)fluoranthene      | 22.875 | 252  | 262666   | 4.515 | ng/ul  | 94       |
| 91) Benzo(k)fluoranthene      | 22.922 | 252  | 272750   | 4.705 | ng/ul  | 98       |
| 93) Benzo(a)pyrene            | 23.457 | 252  | 243658   | 4.628 | ng/ul# | 97       |
| 94) Indeno(1,2,3-cd)pyrene    | 25.827 | 276  | 302548   | 4.194 | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene    | 25.845 | 278  | 257856   | 4.319 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 26.521 | 276  | 266463   | 4.173 | ng/ul# | 95       |

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