

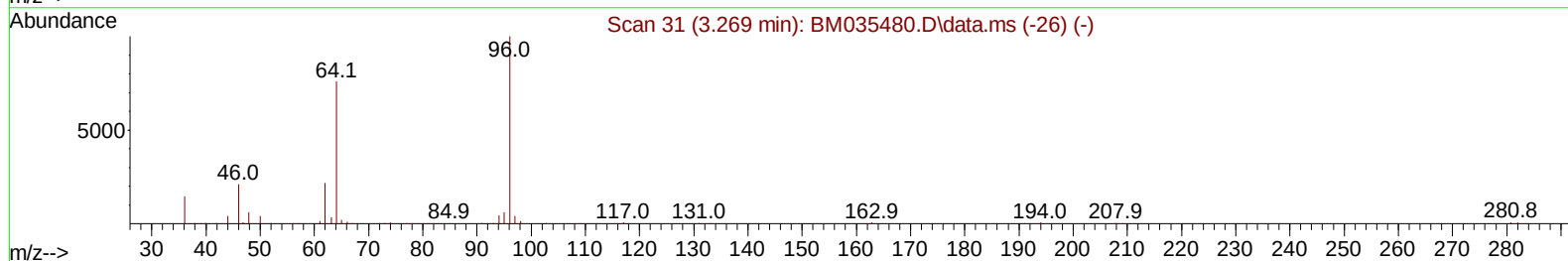
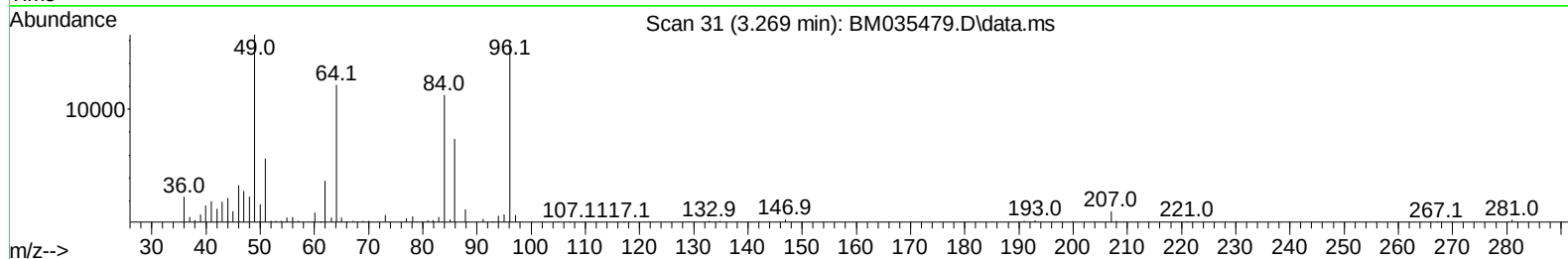
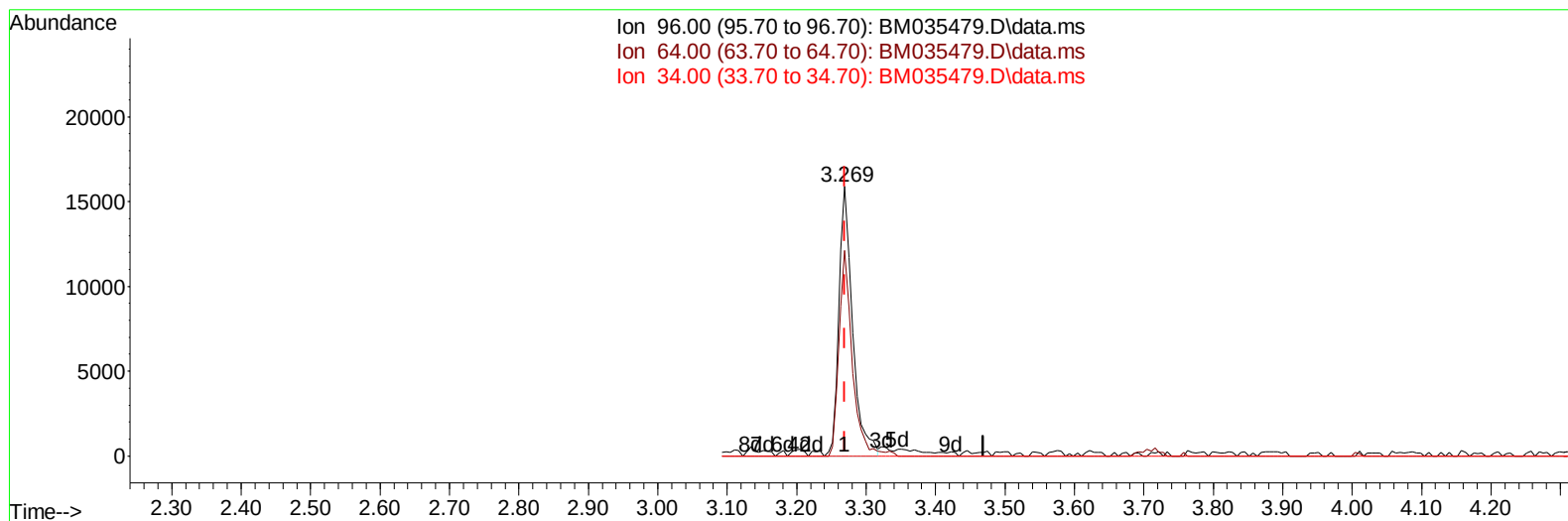
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SST010048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

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

**3.269min (+ 0.000) 3.96 ng/uL**

**response 21614**

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 75.40  | 76.25  |
| 34.00 | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

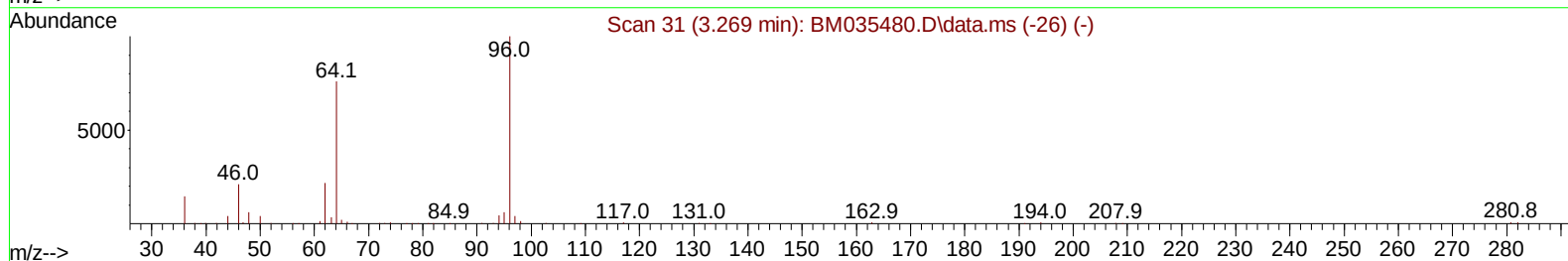
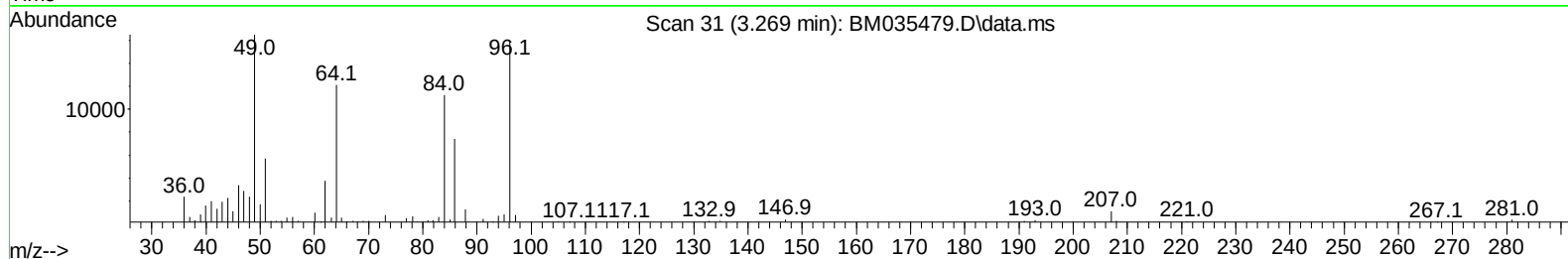
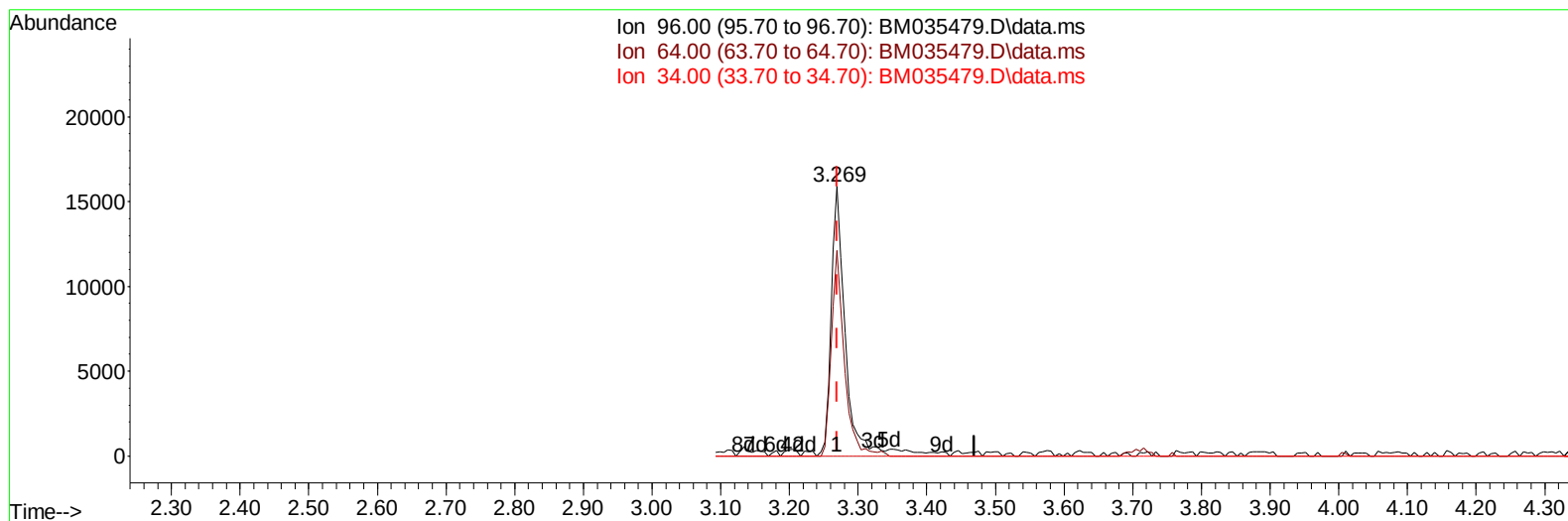
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SST010048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
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Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

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

**3.269min (+ 0.000) 4.07 ng/uL m**

**response 22214**

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 75.40  | 76.25  |
| 34.00 | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

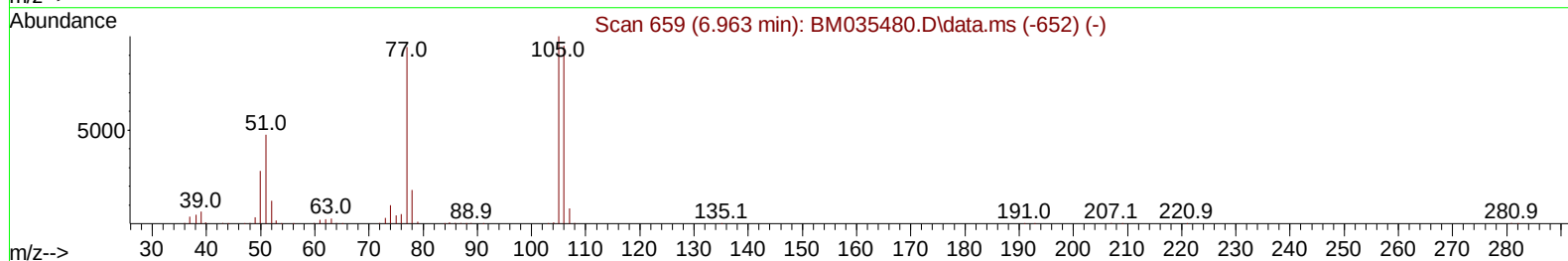
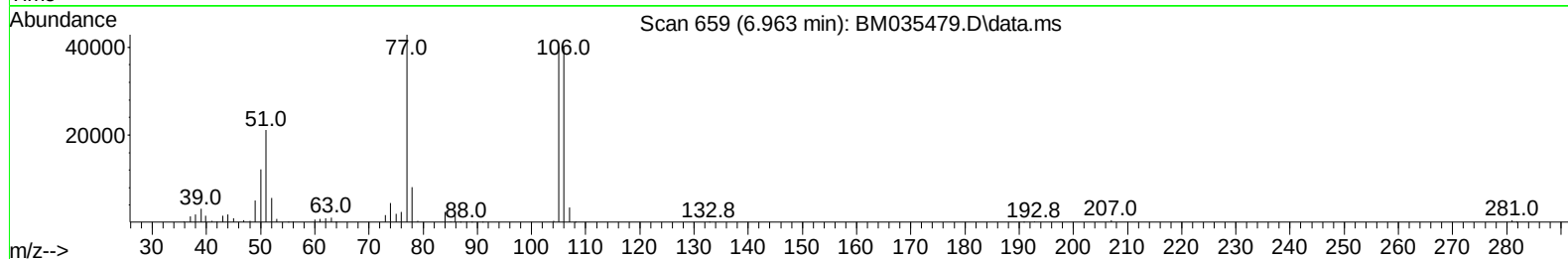
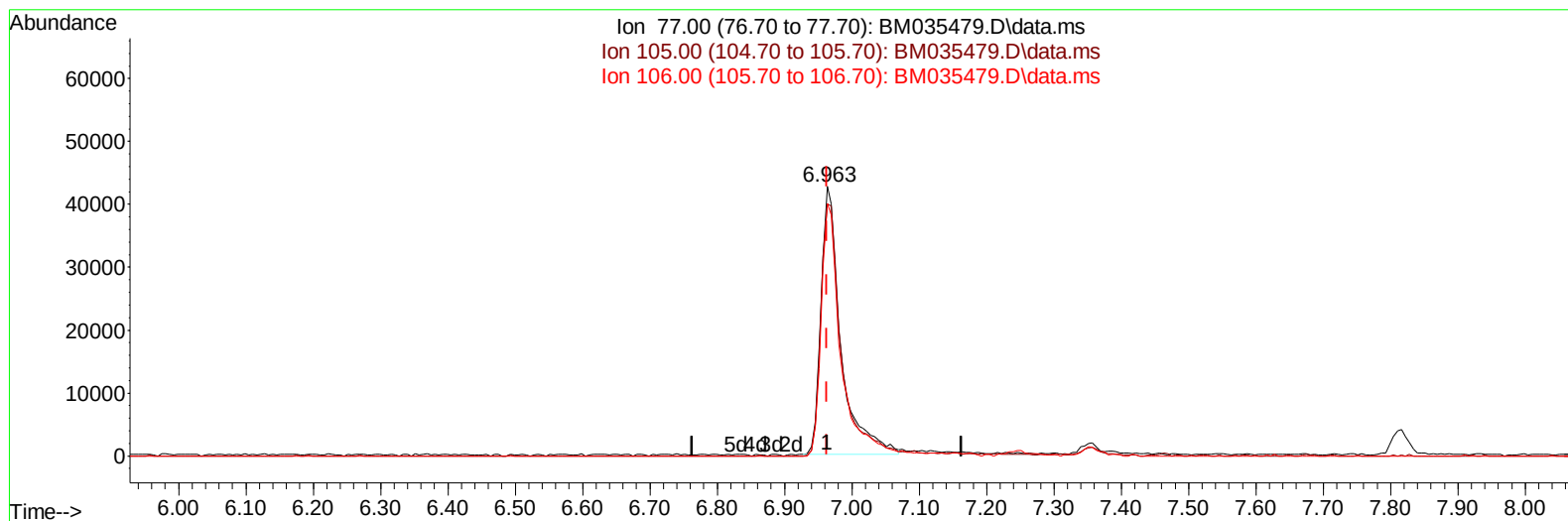
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SSTD01048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

**(6) Benzaldehyde**

**6.963min (+ 0.000) 10.54 ng/u1**

**response 86820**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 105.80 | 93.56  |
| 106.00 | 100.40 | 93.69  |
| 0.00   | 0.00   | 0.00   |

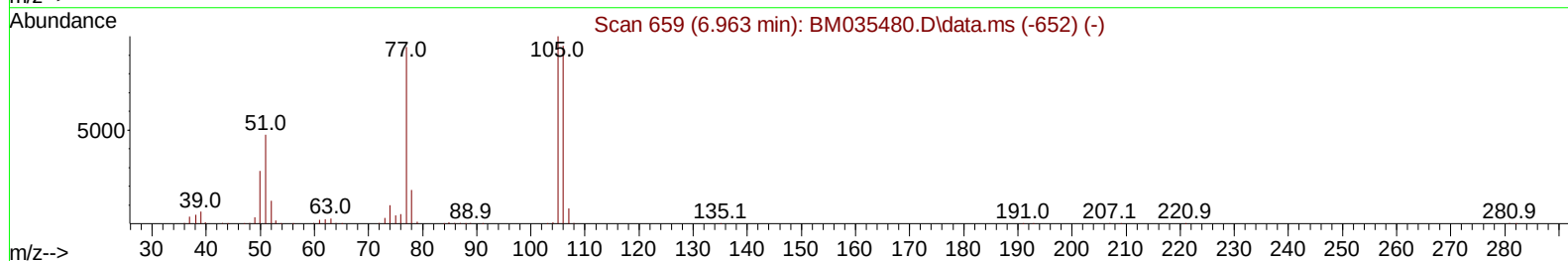
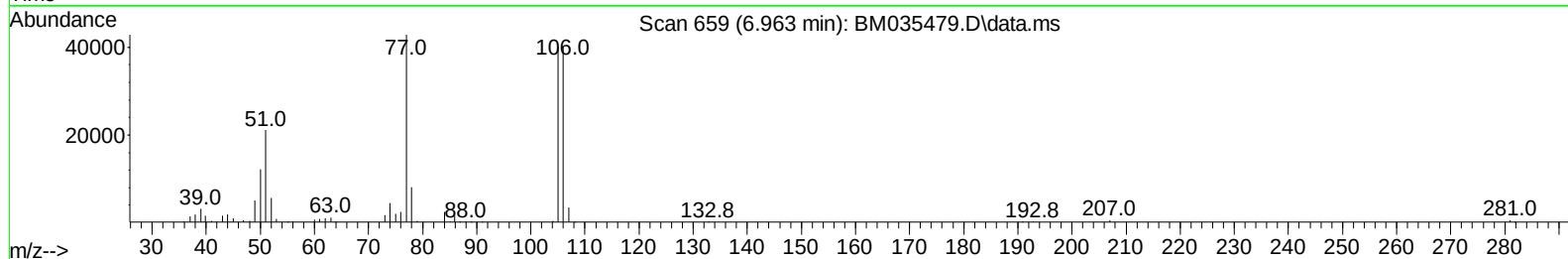
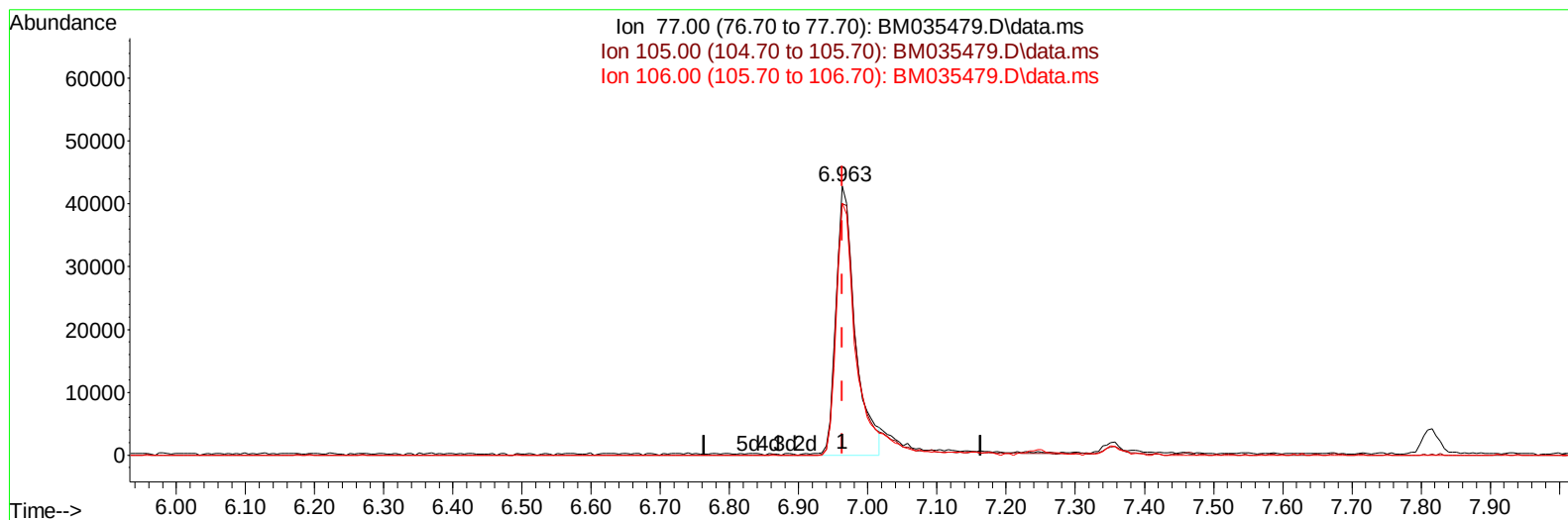
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Data File : BM035479.D  
Acq On : 16 Jun 2022 14:03  
Operator : CG/JU  
Sample : SSTD01048  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Jun 17 00:59:32 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

(6) Benzaldehyde

6.963min (+ 0.000) 9.96 ng/ul m

response 82034

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 105.80 | 93.56  |
| 106.00 | 100.40 | 93.69  |
| 0.00   | 0.00   | 0.00   |

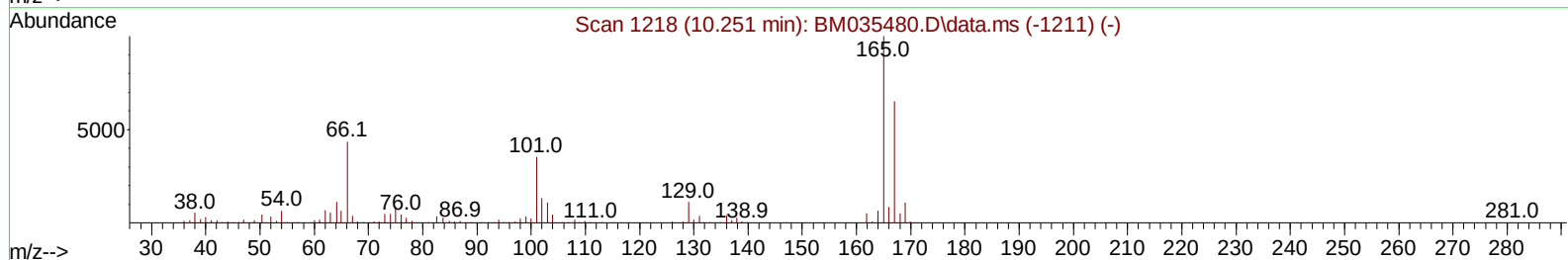
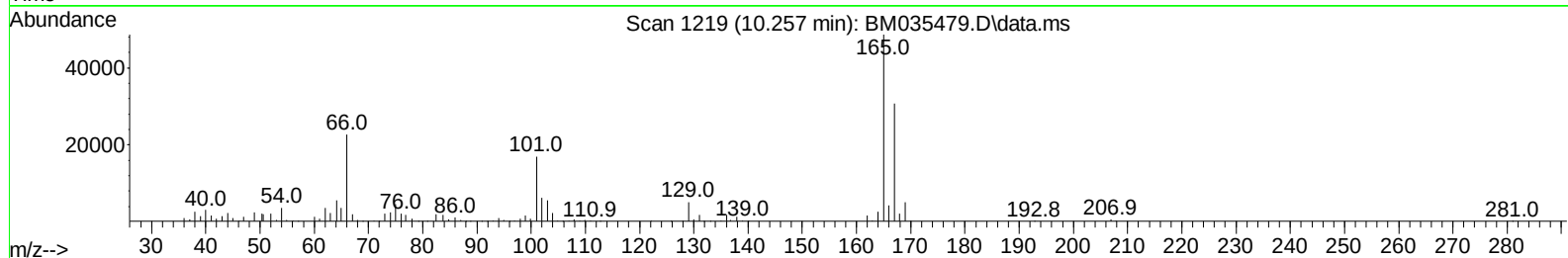
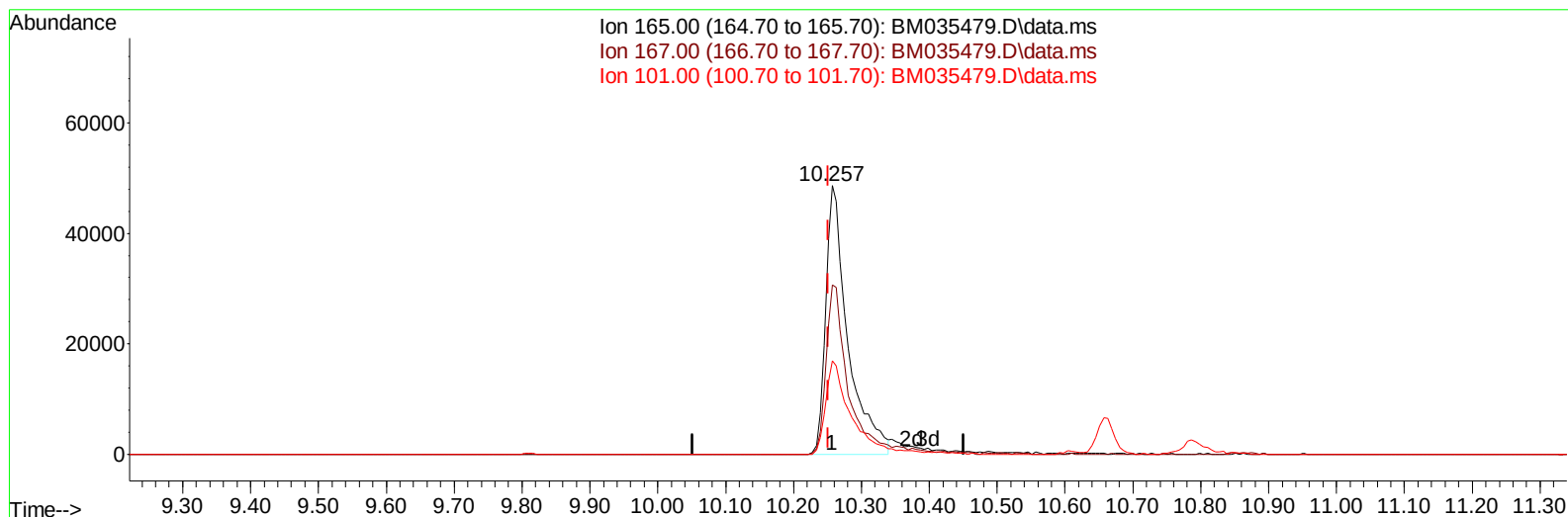
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SSTD01048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.257min (+ 0.006) 7.61 ng/ul

response 110159

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 165.00 | 100.00 | 100.00 |
| 167.00 | 65.20  | 63.10  |
| 101.00 | 35.30  | 34.91  |
| 0.00   | 0.00   | 0.00   |

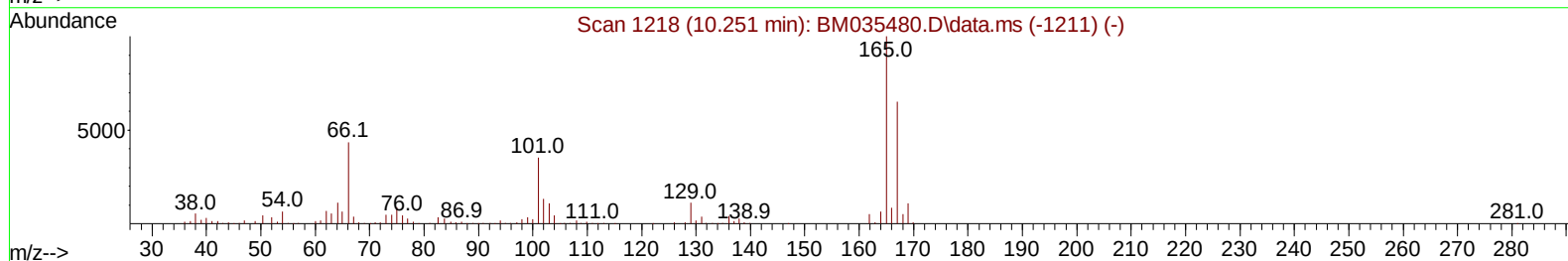
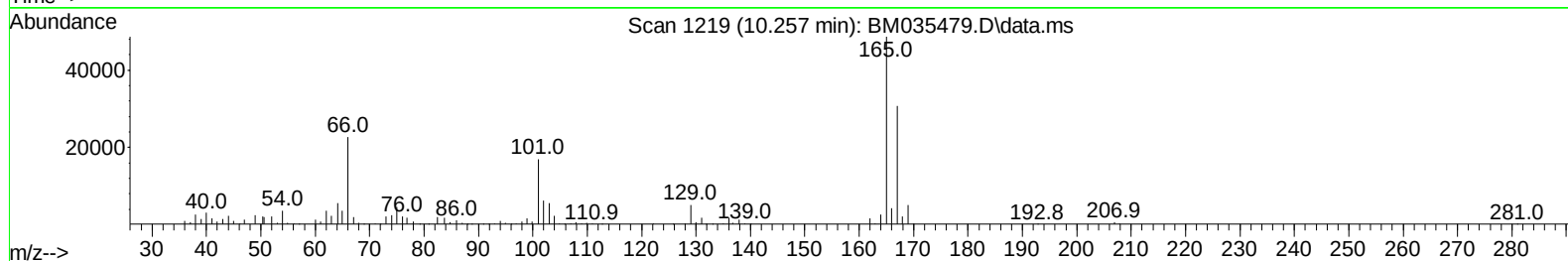
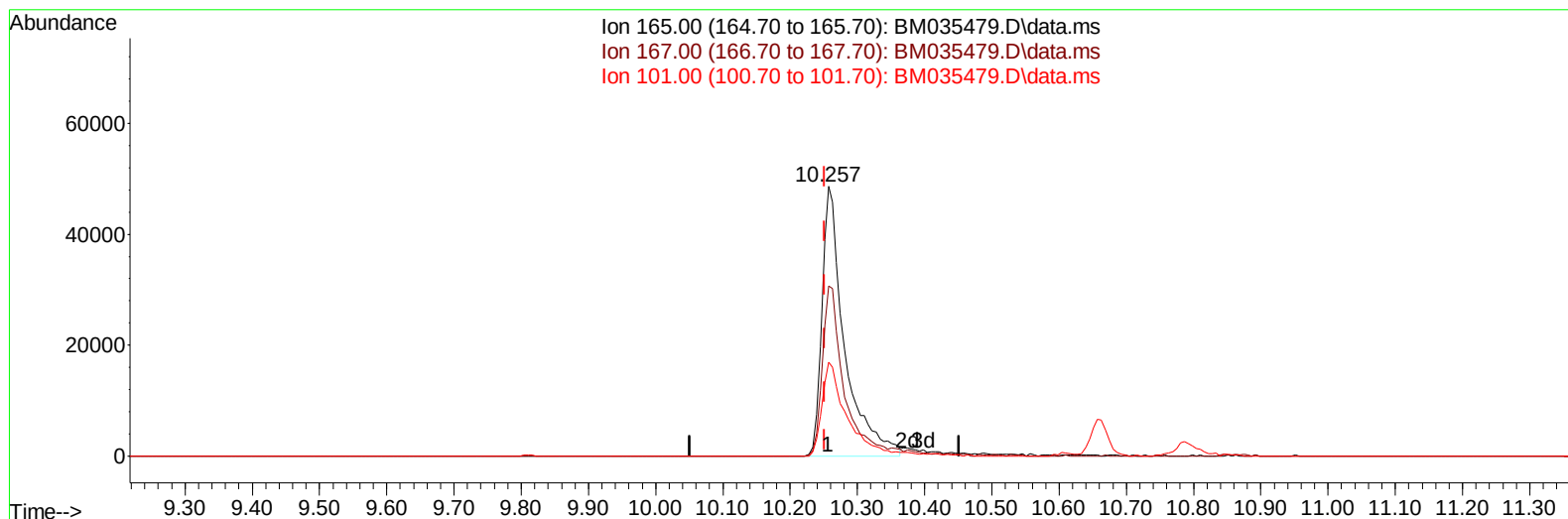
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
Data File : BM035479.D  
Acq On : 16 Jun 2022 14:03  
Operator : CG/JU  
Sample : SSTD01048  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Jun 17 00:59:32 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.257min (+ 0.006) 7.81 ng/ul m

response 113081

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 165.00 | 100.00 | 100.00 |
| 167.00 | 65.20  | 63.10  |
| 101.00 | 35.30  | 34.91  |
| 0.00   | 0.00   | 0.00   |

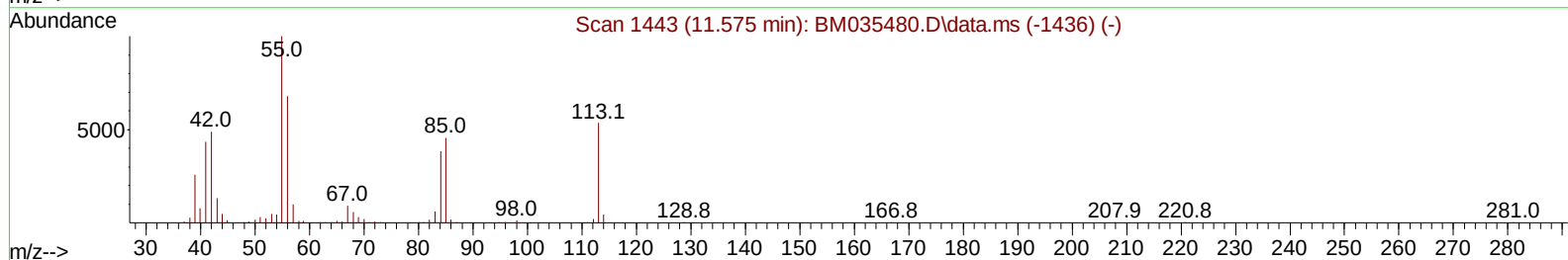
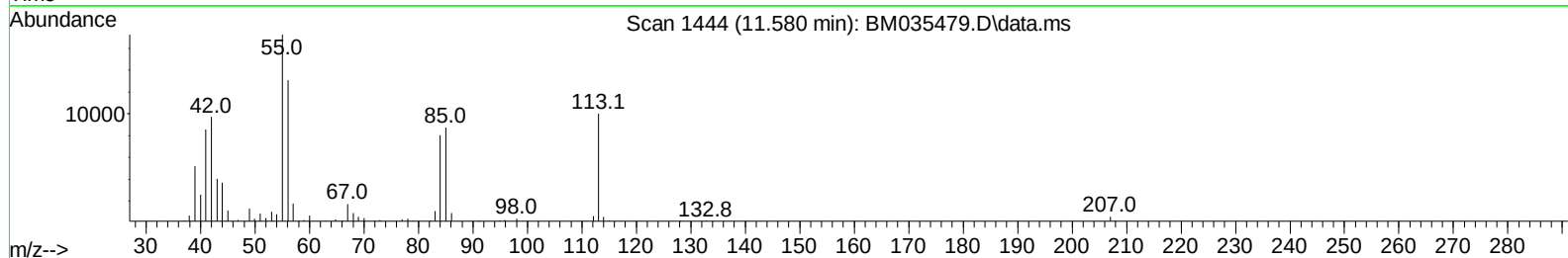
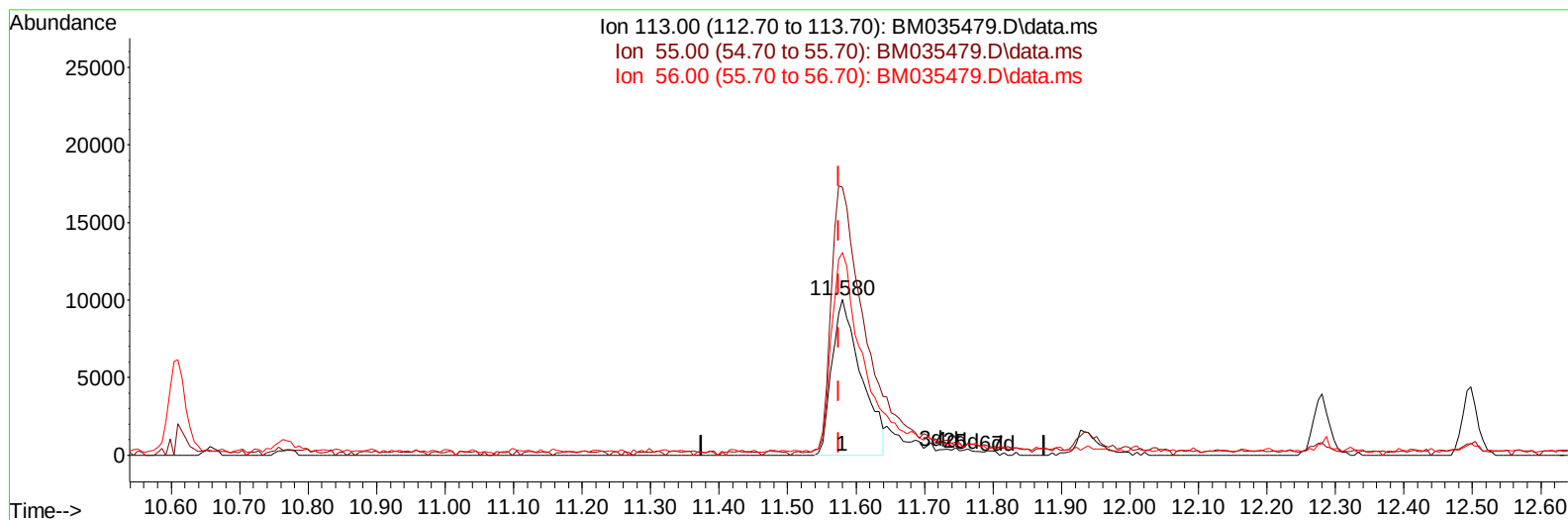
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 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SSTD01048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

**(34) Caprolactam**

**11.580min (+ 0.006) 7.07 ng/ul**

**response 29730**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 187.80 | 172.34 |
| 56.00  | 127.60 | 130.49 |
| 0.00   | 0.00   | 0.00   |

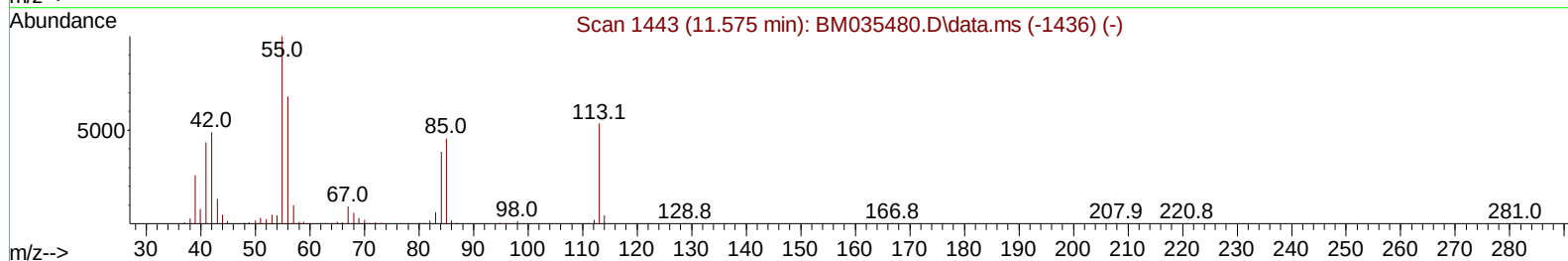
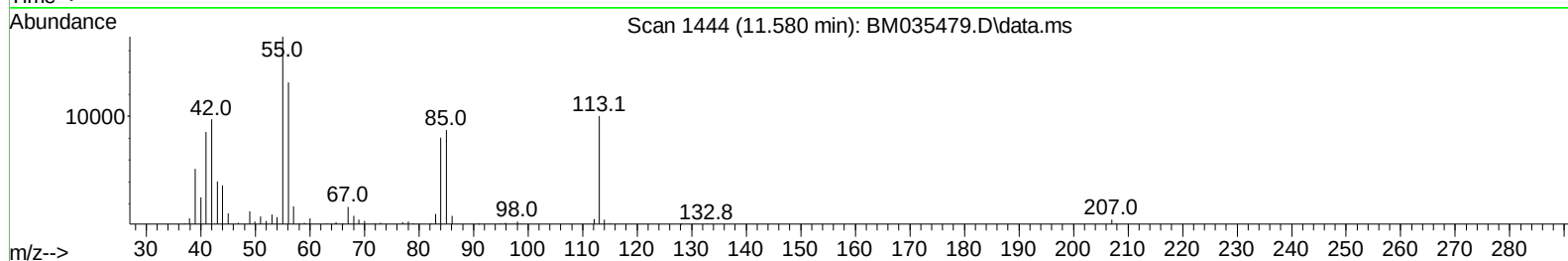
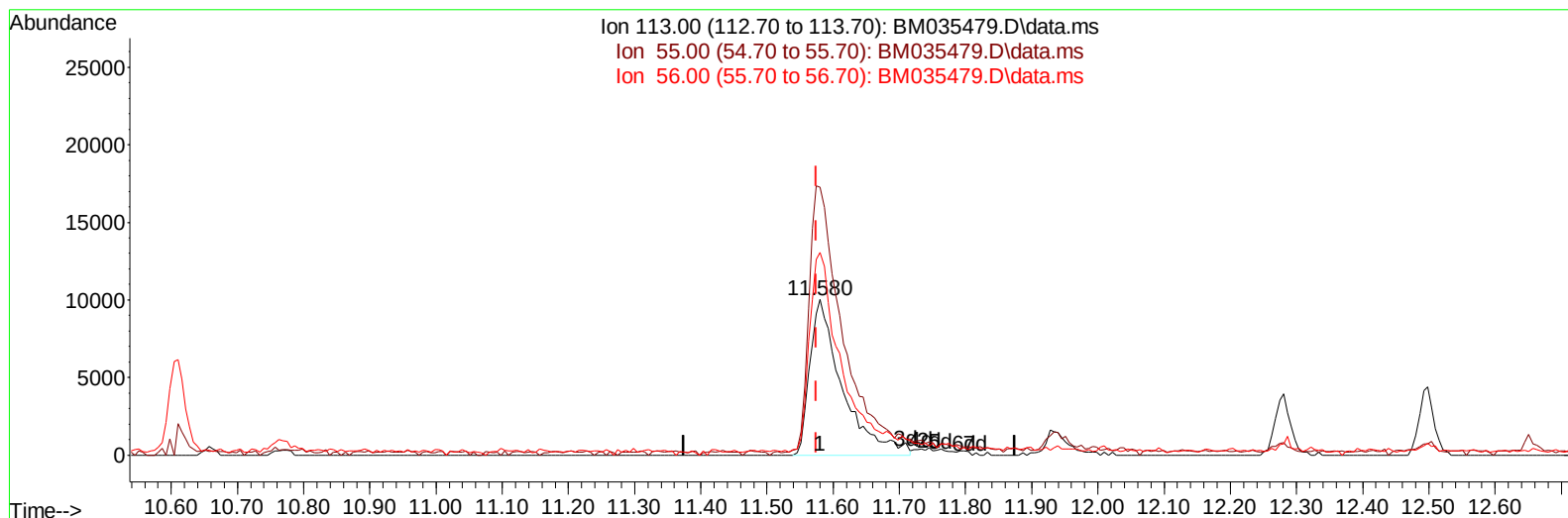
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SSTD01048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

**(34) Caprolactam**

**11.580min (+ 0.006) 8.12 ng/ul m**

**response 34156**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 187.80 | 172.34 |
| 56.00  | 127.60 | 130.49 |
| 0.00   | 0.00   | 0.00   |



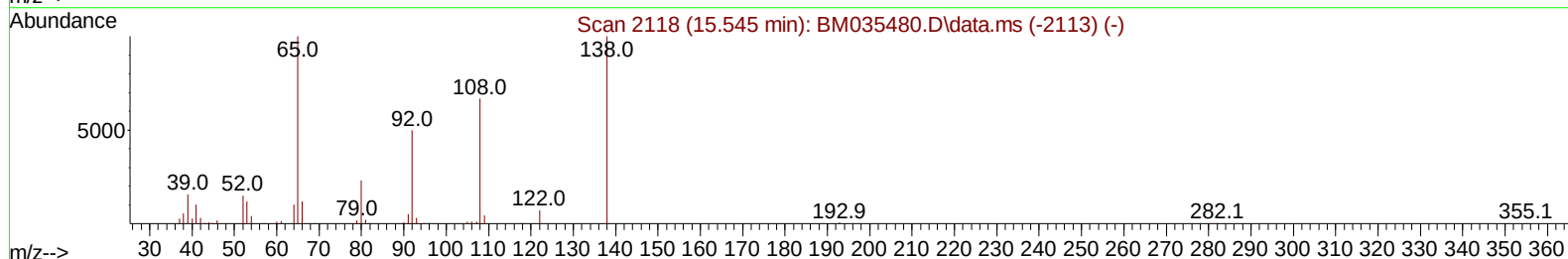
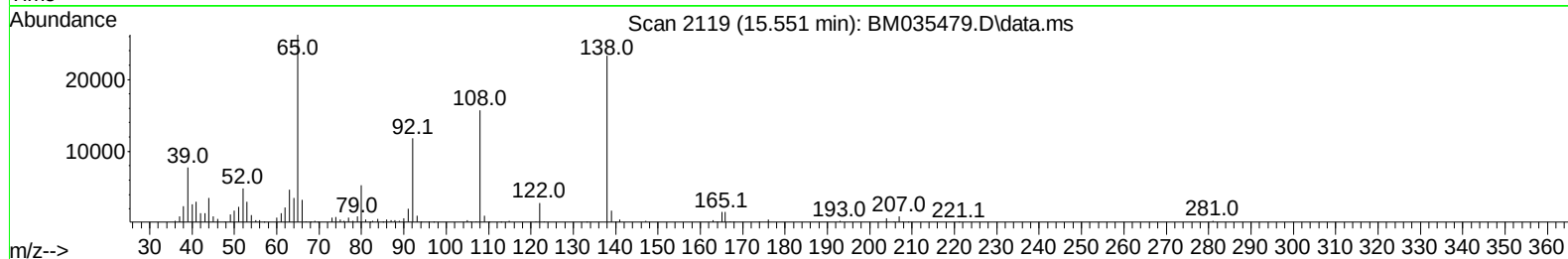
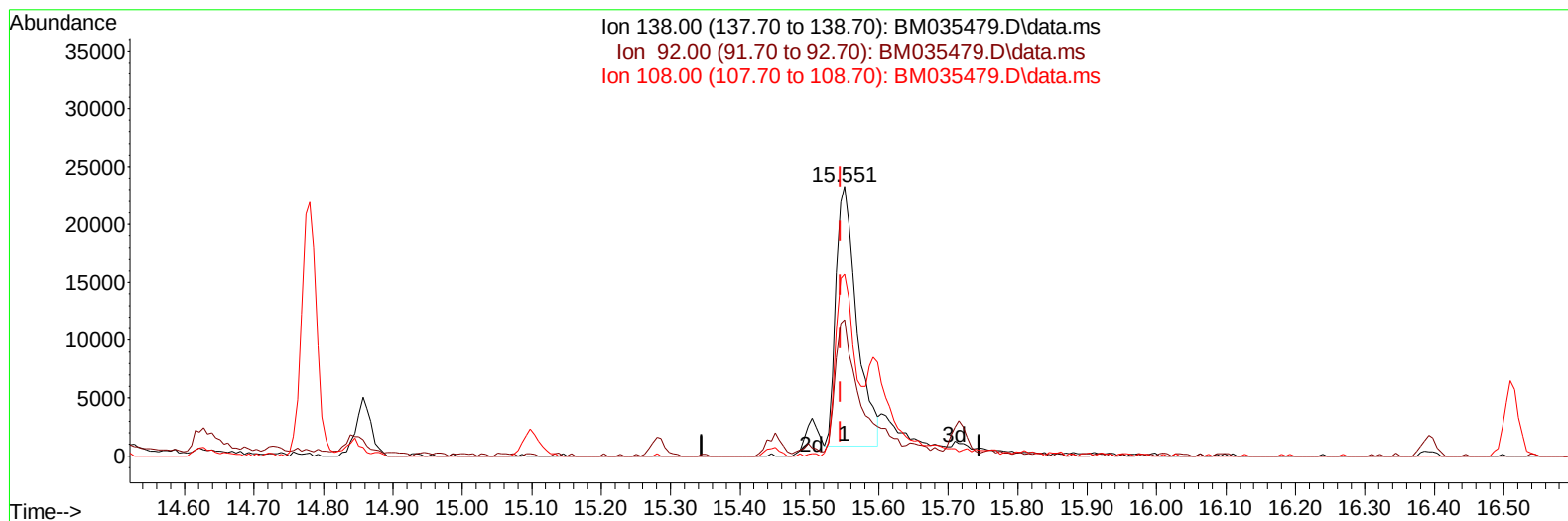
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SSTD01048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

**(63) 4-Nitroaniline**

**15.551min (+ 0.006) 6.73 ng/ul**

**response 46625**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 49.00  | 50.70  |
| 108.00 | 64.50  | 67.63  |
| 0.00   | 0.00   | 0.00   |

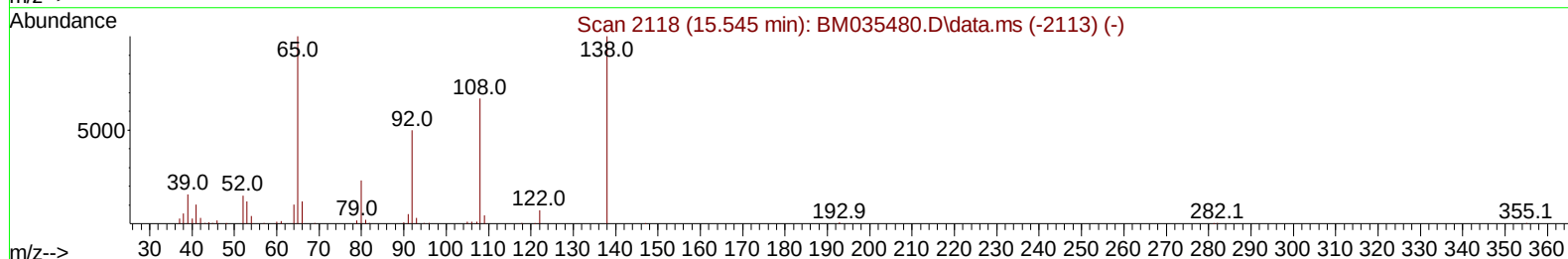
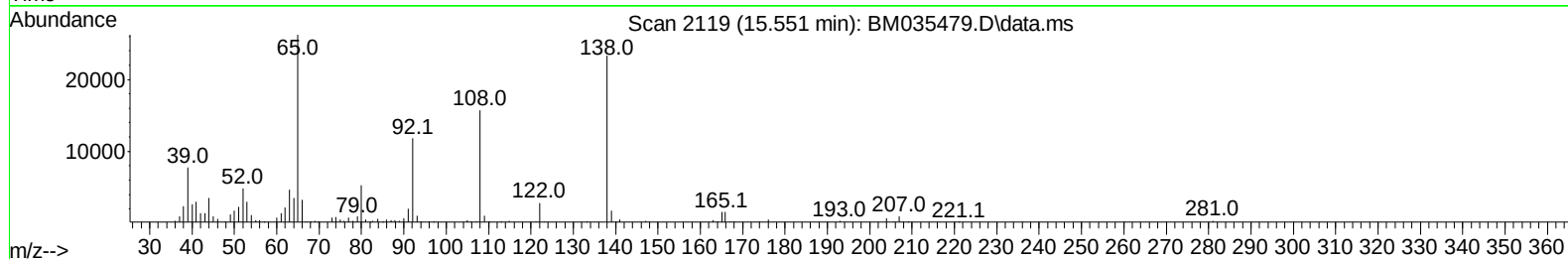
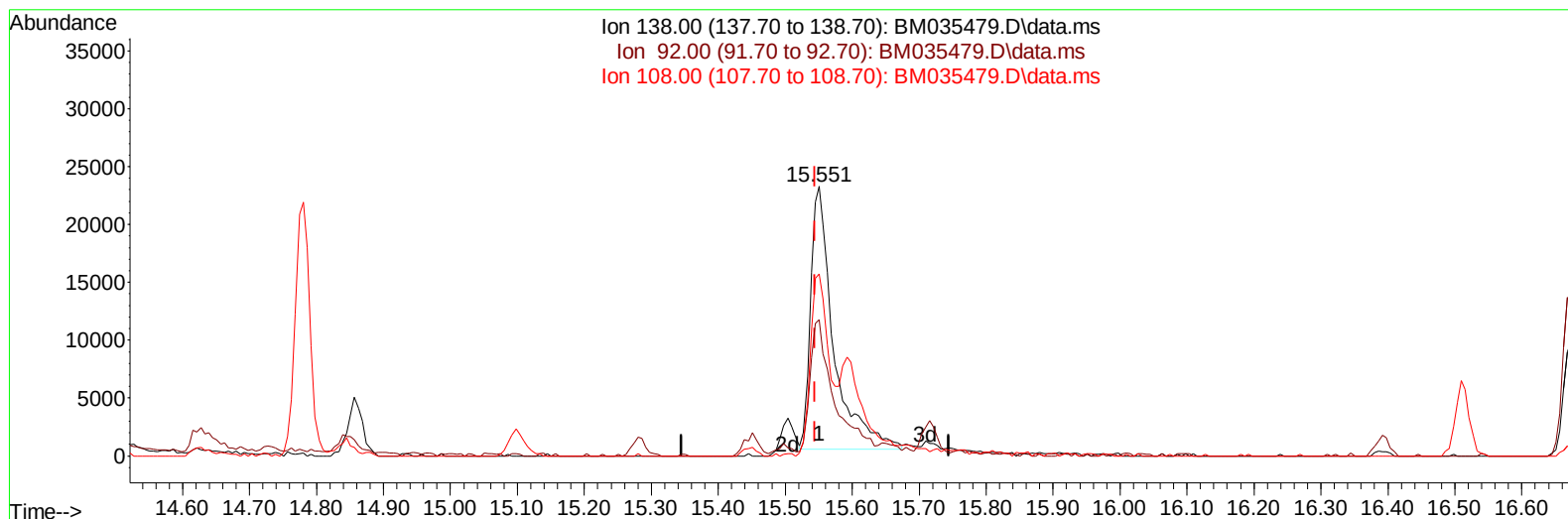
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SSTD01048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022



TIC: BM035479.D\data.ms

**(63) 4-Nitroaniline**

**15.551min (+ 0.006) 7.90 ng/ul m**

**response 54735**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 49.00  | 50.70  |
| 108.00 | 64.50  | 67.63  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SST001048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST0010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.816  | 152  | 225257   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.610 | 136  | 927537   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 548923   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.204 | 188  | 1108522  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.392 | 240  | 1156002  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 23.733 | 264  | 1247951  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.269  | 96   | 22214m   | 4.075  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.693  | 84   | 108380   | 7.707  | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.004  | 99   | 147879   | 8.908  | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.151  | 67   | 95214    | 9.090  | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.351  | 132  | 136441   | 9.271  | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.545  | 113  | 121986   | 8.747  | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.981  | 128  | 62422    | 8.584  | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.704  | 143  | 64555    | 7.928  | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.257 | 165  | 113081m  | 7.808  | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.763 | 131  | 175656   | 8.796  | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.869 | 166  | 351041   | 8.778  | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.151 | 160  | 432479   | 9.048  | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.721 | 143  | 28373    | 4.002  | ng/ul  | 0.02     |
| 60) Fluorene-d10              | 15.451 | 176  | 311934   | 9.220  | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.592 | 200  | 47783    | 6.675  | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.304 | 188  | 472419   | 9.156  | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.598 | 212  | 534171   | 8.676  | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.586 | 264  | 564148   | 8.920  | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.305  | 88   | 20806    | 3.673  | ng/uL  | 92       |
| 5) Pyridine                   | 3.710  | 79   | 123201   | 8.452  | ng/ul  | 92       |
| 6) Benzaldehyde               | 6.963  | 77   | 82034m   | 9.959  | ng/ul  |          |
| 8) Phenol                     | 7.034  | 94   | 156960   | 8.990  | ng/ul  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.240  | 93   | 129901   | 9.599  | ng/ul  | 98       |
| 12) 2-Chlorophenol            | 7.387  | 128  | 140176   | 9.323  | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.275  | 108  | 118874   | 9.071  | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.351  | 45   | 200995   | 10.455 | ng/ul  | 99       |
| 16) Acetophenone              | 8.645  | 105  | 196467   | 9.095  | ng/ul  | 94       |
| 17) N-Nitroso-di-n-propyla... | 8.628  | 70   | 92115    | 8.995  | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.610  | 108  | 125995   | 8.997  | ng/ul  | 96       |
| 19) Hexachloroethane          | 8.887  | 117  | 57900    | 8.859  | ng/ul  | 91       |
| 22) Nitrobenzene              | 9.022  | 77   | 132868   | 7.641  | ng/ul  | 97       |
| 23) Isophorone                | 9.545  | 82   | 249249   | 8.121  | ng/ul  | 100      |
| 25) 2-Nitrophenol             | 9.734  | 139  | 71961    | 8.255  | ng/ul  | 98       |
| 26) 2,4-Dimethylphenol        | 9.804  | 107  | 140641   | 8.186  | ng/ul  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 10.028 | 93   | 164911   | 8.927  | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.286 | 162  | 118696   | 8.128  | ng/ul  | 98       |
| 30) Naphthalene               | 10.657 | 128  | 449613   | 9.372  | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.786 | 127  | 173265   | 8.676  | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 10.945 | 225  | 77298    | 7.279  | ng/ul  | 98       |
| 34) Caprolactam               | 11.580 | 113  | 34156m   | 8.118  | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.933 | 107  | 118799   | 8.103  | ng/ul  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061622\  
 Data File : BM035479.D  
 Acq On : 16 Jun 2022 14:03  
 Operator : CG/JU  
 Sample : SST001048  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST0010048

Manual IntegrationsAPPROVED

Quant Time: Jun 17 01:04:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 00:59:32 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/21/2022  
 Supervised By :Yogesh Patel 06/22/2022

| Compound                      | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.280 | 142  | 291501   | 9.070 | ng/ul  | 98       |
| 37) 1-Methylnaphthalene       | 12.498 | 142  | 296490   | 9.204 | ng/ul  | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.651 | 216  | 143470   | 8.128 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.627 | 237  | 26664    | 2.118 | ng/ul  | 92       |
| 41) 2,4,6-Trichlorophenol     | 12.910 | 196  | 84897    | 7.531 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol     | 12.998 | 196  | 88023    | 7.357 | ng/ul  | 97       |
| 43) 1,1'-Biphenyl             | 13.292 | 154  | 384638   | 9.051 | ng/ul  | 100      |
| 44) 2-Chloronaphthalene       | 13.333 | 162  | 297392   | 9.024 | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.557 | 65   | 67094    | 7.246 | ng/ul  | 96       |
| 47) Dimethylphthalate         | 13.921 | 163  | 359822   | 8.986 | ng/ul  | 100      |
| 48) 2,6-Dinitrotoluene        | 14.045 | 165  | 64724    | 8.115 | ng/ul  | 95       |
| 50) Acenaphthylene            | 14.174 | 152  | 484437   | 9.431 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.386 | 138  | 60115    | 7.860 | ng/ul  | 97       |
| 52) Acenaphthene              | 14.521 | 153  | 317407   | 9.278 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.627 | 184  | 21751    | 4.163 | ng/ul  | 97       |
| 55) 4-Nitrophenol             | 14.745 | 109  | 15055    | 2.264 | ng/ul  | 93       |
| 56) Dibenzofuran              | 14.857 | 168  | 447647   | 9.356 | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene        | 14.845 | 165  | 97034    | 8.664 | ng/ul# | 87       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.098 | 232  | 76672    | 7.910 | ng/ul# | 97       |
| 59) Diethylphthalate          | 15.280 | 149  | 364858   | 8.998 | ng/ul  | 99       |
| 61) Fluorene                  | 15.504 | 166  | 366187   | 9.407 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.498 | 204  | 172481   | 8.728 | ng/ul  | 99       |
| 63) 4-Nitroaniline            | 15.551 | 138  | 54735m   | 7.899 | ng/ul  |          |
| 66) 4,6-Dinitro-2-methylph... | 15.610 | 198  | 48832    | 6.806 | ng/ul# | 98       |
| 67) N-Nitrosodiphenylamine    | 15.715 | 169  | 309095   | 9.217 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.392 | 248  | 99006    | 8.433 | ng/ul  | 99       |
| 69) Hexachlorobenzene         | 16.515 | 284  | 117655   | 8.776 | ng/ul  | 98       |
| 70) Atrazine                  | 16.674 | 200  | 102713   | 7.850 | ng/ul  | 97       |
| 71) Pentachlorophenol         | 16.880 | 266  | 48019    | 5.360 | ng/ul  | 95       |
| 72) Phenanthrene              | 17.245 | 178  | 575108   | 9.581 | ng/ul  | 99       |
| 74) Anthracene                | 17.339 | 178  | 582912   | 9.530 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.257 | 216  | 158329   | 8.199 | ng/uL  | 98       |
| 76) Pentachlorobenzene        | 14.780 | 250  | 143006   | 8.621 | ng/uL  | 97       |
| 77) Carbazole                 | 17.615 | 167  | 515449   | 9.522 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.174 | 149  | 658688   | 8.503 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.268 | 202  | 648571   | 8.770 | ng/ul  | 99       |
| 82) Pyrene                    | 19.627 | 202  | 680630   | 8.941 | ng/ul  | 99       |
| 83) Butylbenzylphthalate      | 20.527 | 149  | 279178   | 8.562 | ng/ul  | 96       |
| 84) 3,3'-Dichlorobenzidine    | 21.315 | 252  | 208613   | 8.778 | ng/ul  | 98       |
| 85) Benzo(a)anthracene        | 21.380 | 228  | 660270   | 8.940 | ng/ul  | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.309 | 149  | 421418   | 8.595 | ng/ul  | 99       |
| 87) Chrysene                  | 21.433 | 228  | 668849   | 9.289 | ng/ul  | 100      |
| 89) Di-n-octyl phthalate      | 22.209 | 149  | 739499   | 8.265 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.027 | 252  | 728451   | 8.975 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.074 | 252  | 661648   | 8.581 | ng/ul  | 98       |
| 93) Benzo(a)pyrene            | 23.633 | 252  | 699419   | 8.845 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.156 | 276  | 823050   | 9.058 | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene    | 26.162 | 278  | 704715   | 9.029 | ng/ul  | 99       |
| 96) Benzo(g,h,i)perylene      | 26.897 | 276  | 697215   | 9.224 | ng/ul  | 100      |

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

**Instrument :**

BNA\_M

**ClientSampleId :**

SSTD010048

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 06/21/2022  
Supervised By :Yogesh Patel 06/22/2022

