

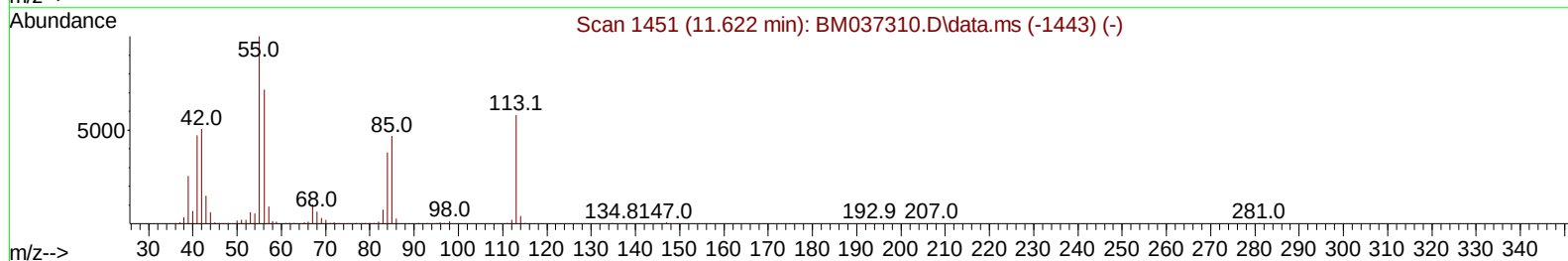
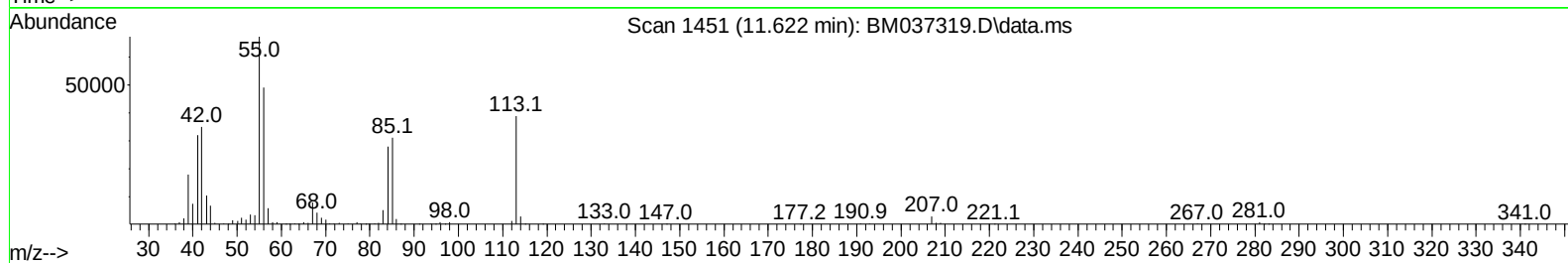
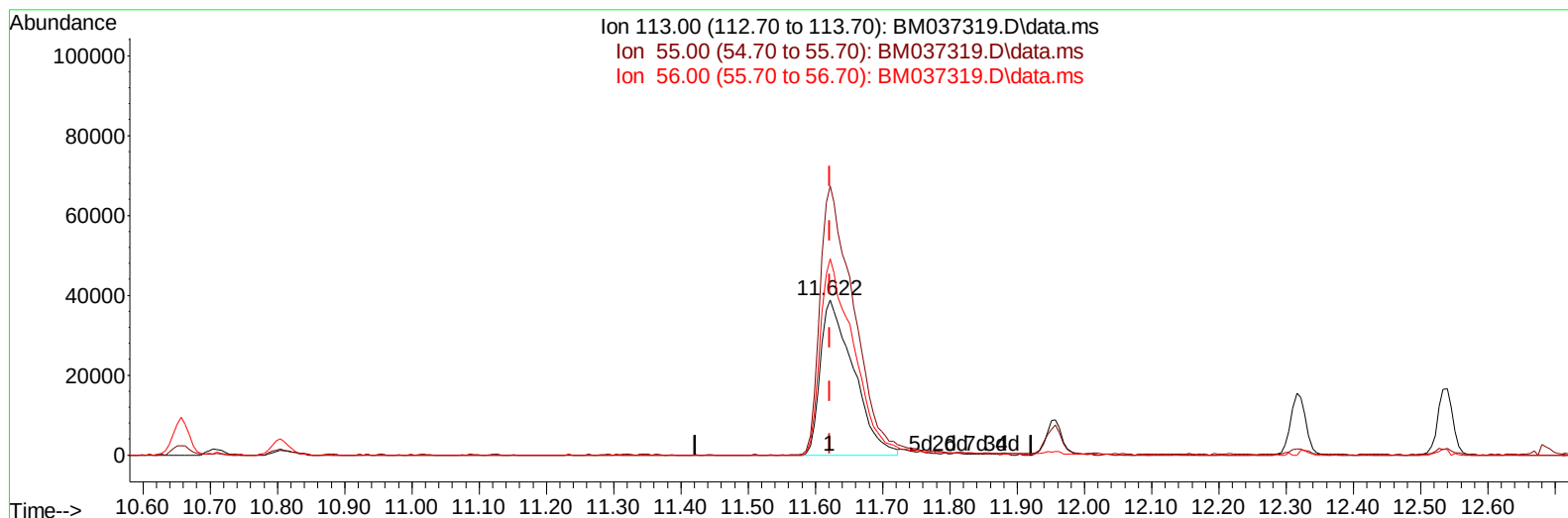
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110222\  
 Data File : BM037319.D  
 Acq On : 02 Nov 2022 16:53  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Nov 02 22:54:19 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 02 03:33:37 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2022  
 Supervised By :mohammad ahmed 11/04/2022



TIC: BM037319.D\data.ms

**(34) Caprolactam**

**11.622min (+ 0.000) 18.84 ng/ul**

**response 131960**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 172.10 | 173.22 |
| 56.00  | 121.30 | 126.36 |
| 0.00   | 0.00   | 0.00   |

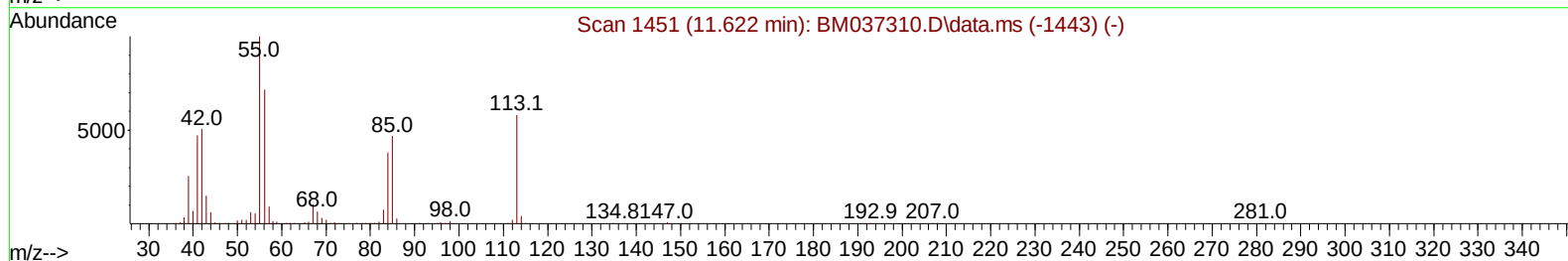
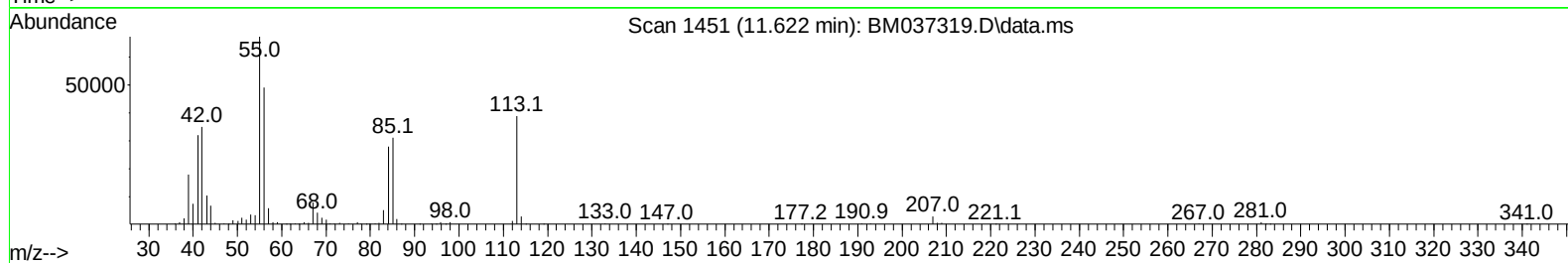
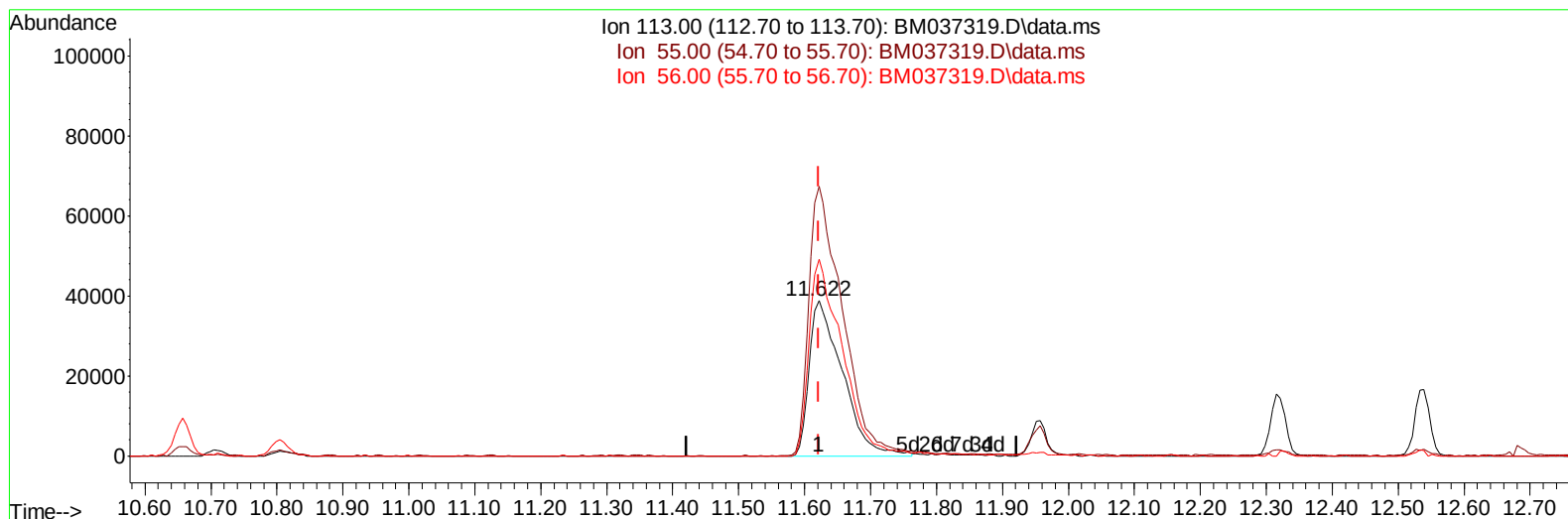
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110222\  
 Data File : BM037319.D  
 Acq On : 02 Nov 2022 16:53  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Nov 02 22:54:19 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 02 03:33:37 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2022  
 Supervised By :mohammad ahmed 11/04/2022



TIC: BM037319.D\data.ms

**(34) Caprolactam**

11.622min (+ 0.000) 19.19 ng/ul m

response 134448

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 172.10 | 173.22 |
| 56.00  | 121.30 | 126.36 |
| 0.00   | 0.00   | 0.00   |

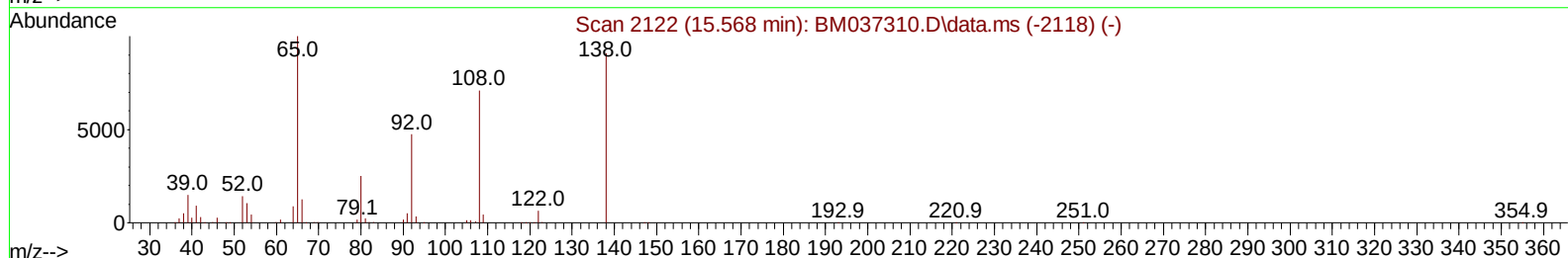
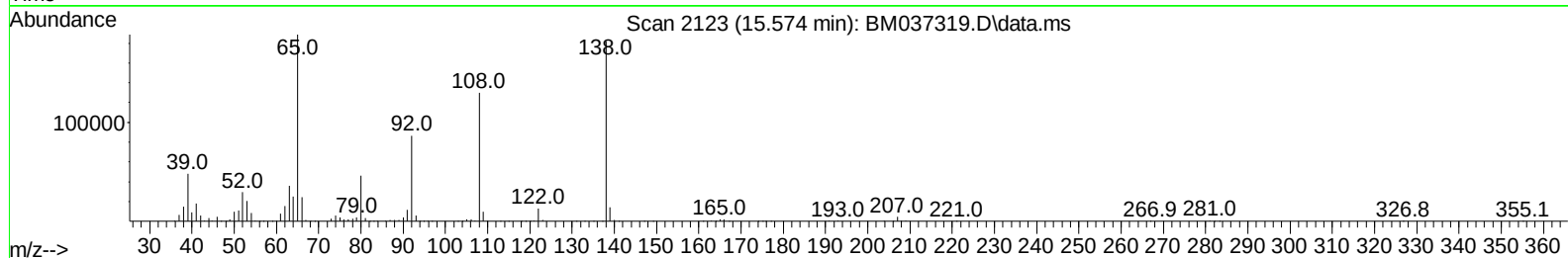
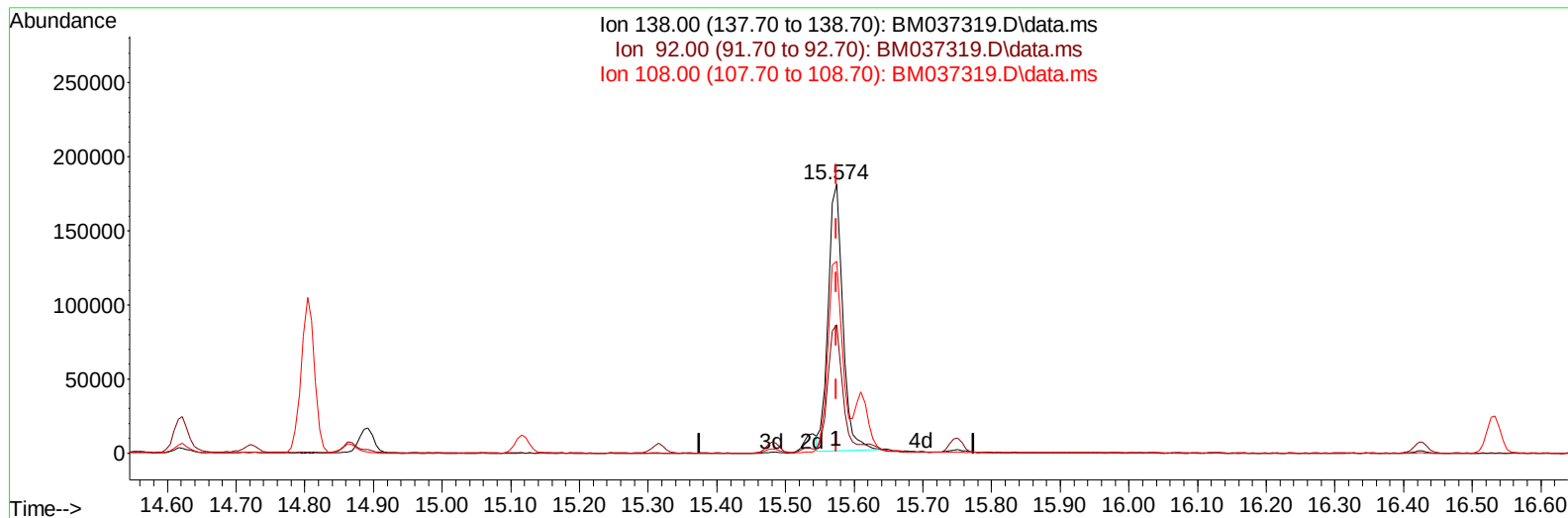
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110222\  
Data File : BM037319.D  
Acq On : 02 Nov 2022 16:53  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_M  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Nov 02 22:54:19 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110122.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Nov 02 03:33:37 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2022  
Supervised By :mohammad ahmed 11/04/2022



TIC: BM037319.D\data.ms

**(63) 4-Nitroaniline**

**15.574min (+ 0.000) 20.84 ng/ul**

**response 264859**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 48.10  | 47.69  |
| 108.00 | 71.20  | 71.41  |
| 0.00   | 0.00   | 0.00   |

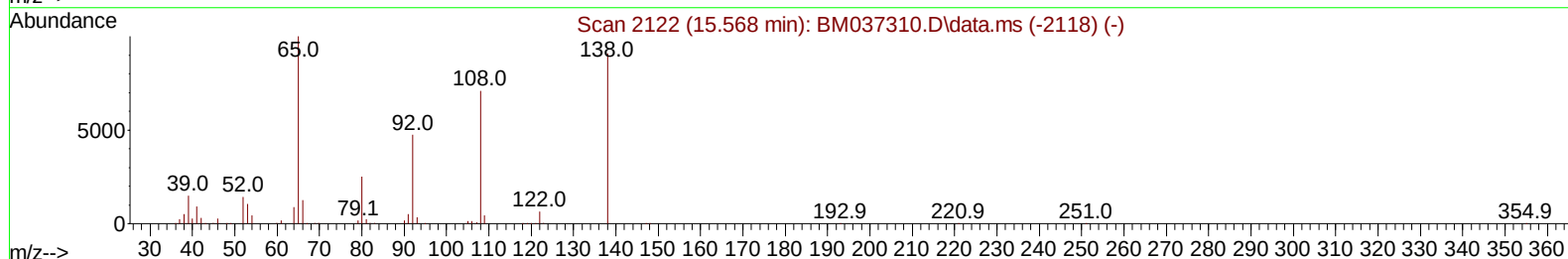
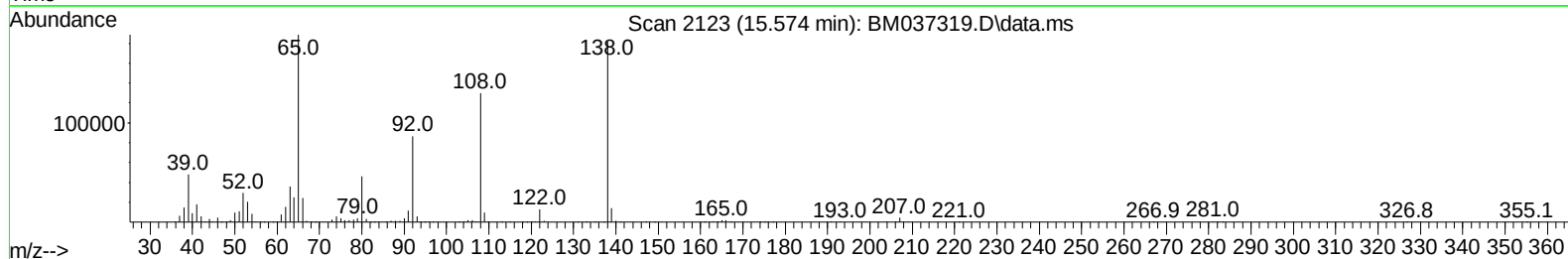
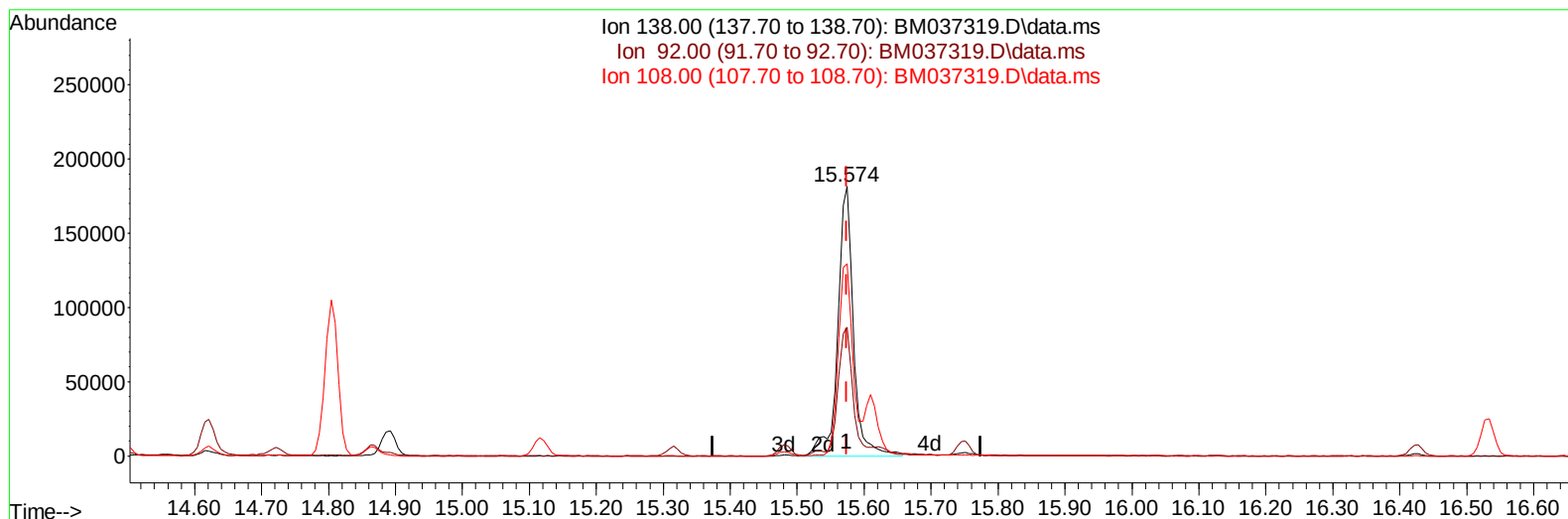
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110222\  
 Data File : BM037319.D  
 Acq On : 02 Nov 2022 16:53  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Nov 02 22:54:19 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 02 03:33:37 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2022  
 Supervised By :mohammad ahmed 11/04/2022



TIC: BM037319.D\data.ms

**(63) 4-Nitroaniline**

**15.574min (+ 0.000) 23.22 ng/ul m**

**response 295118**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 48.10  | 47.69  |
| 108.00 | 71.20  | 71.41  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110222\  
 Data File : BM037319.D  
 Acq On : 02 Nov 2022 16:53  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Nov 02 22:54:19 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 02 03:33:37 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2022  
 Supervised By :mohammad ahmed 11/04/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.857  | 152  | 299065   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.657 | 136  | 1378880  | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.492 | 164  | 878619   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.233 | 188  | 1811853  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.403 | 240  | 1733297  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 23.756 | 264  | 1399880  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.269  | 96   | 57117    | 8.419  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.699  | 84   | 432008   | 21.332 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.028  | 99   | 510028   | 19.596 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.193  | 67   | 329032   | 20.602 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.387  | 132  | 415246   | 20.882 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.575  | 113  | 419911   | 19.757 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.028  | 128  | 214190   | 20.795 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.745  | 143  | 227632   | 20.289 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.281 | 165  | 432670   | 20.917 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.804 | 131  | 639068   | 20.217 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.910 | 166  | 1268869  | 20.887 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.186 | 160  | 1558063  | 21.333 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.704 | 143  | 237300   | 19.542 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.480 | 176  | 1159269  | 21.399 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.610 | 200  | 214562   | 18.697 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.333 | 188  | 1781373  | 21.348 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.621 | 212  | 1950417  | 21.534 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.603 | 264  | 1539117  | 21.359 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.310  | 88   | 62209    | 8.728  | ng/uL  | 99       |
| 5) Pyridine                   | 3.716  | 79   | 424577   | 20.931 | ng/ul  | 97       |
| 6) Benzaldehyde               | 7.004  | 77   | 314586   | 26.127 | ng/ul  | 97       |
| 8) Phenol                     | 7.057  | 94   | 539112   | 19.687 | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.287  | 93   | 449605   | 20.465 | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.416  | 128  | 453864   | 21.093 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.304  | 108  | 413133   | 19.774 | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.392  | 45   | 651774   | 20.889 | ng/ul  | 99       |
| 16) Acetophenone              | 8.687  | 105  | 699020   | 20.508 | ng/ul  | 97       |
| 17) N-Nitroso-di-n-propyla... | 8.675  | 70   | 354508   | 20.939 | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.640  | 108  | 461853   | 20.068 | ng/ul  | 100      |
| 19) Hexachloroethane          | 8.928  | 117  | 186364   | 21.823 | ng/ul  | 96       |
| 22) Nitrobenzene              | 9.069  | 77   | 530611   | 21.344 | ng/ul  | 99       |
| 23) Isophorone                | 9.592  | 82   | 981830   | 20.477 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.775  | 139  | 266317   | 20.768 | ng/ul  | 98       |
| 26) 2,4-Dimethylphenol        | 9.839  | 107  | 512672   | 21.262 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 10.081 | 93   | 629231   | 21.108 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.310 | 162  | 447109   | 21.004 | ng/ul  | 99       |
| 30) Naphthalene               | 10.710 | 128  | 1560581  | 21.213 | ng/ul  | 100      |
| 32) 4-Chloroaniline           | 10.828 | 127  | 668636   | 20.413 | ng/ul  | 99       |
| 33) Hexachlorobutadiene       | 10.986 | 225  | 280145   | 21.370 | ng/ul  | 99       |
| 34) Caprolactam               | 11.622 | 113  | 134448m  | 19.193 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.957 | 107  | 465393   | 21.078 | ng/ul  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110222\  
 Data File : BM037319.D  
 Acq On : 02 Nov 2022 16:53  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Nov 02 22:54:19 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 02 03:33:37 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2022  
 Supervised By :mohammad ahmed 11/04/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.316 | 142  | 1068121  | 21.024 | ng/ul | 99       |
| 37) 1-Methylnaphthalene       | 12.539 | 142  | 1075070  | 21.029 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.686 | 216  | 553378   | 20.985 | ng/ul | 100      |
| 40) Hexachlorocyclopentadiene | 12.657 | 237  | 279416   | 18.953 | ng/ul | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.933 | 196  | 356374   | 20.935 | ng/ul | 99       |
| 42) 2,4,5-Trichlorophenol     | 13.004 | 196  | 387832   | 21.159 | ng/ul | 99       |
| 43) 1,1'-Biphenyl             | 13.327 | 154  | 1403989  | 21.500 | ng/ul | 100      |
| 44) 2-Chloronaphthalene       | 13.369 | 162  | 1118150  | 21.309 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.586 | 65   | 306952   | 22.185 | ng/ul | 98       |
| 47) Dimethylphthalate         | 13.957 | 163  | 1366698  | 21.121 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 14.080 | 165  | 268111   | 21.161 | ng/ul | 97       |
| 50) Acenaphthylene            | 14.216 | 152  | 1712196  | 21.449 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.410 | 138  | 293478   | 21.464 | ng/ul | 98       |
| 52) Acenaphthene              | 14.557 | 153  | 1211908  | 21.336 | ng/ul | 99       |
| 53) 2,4-Dinitrophenol         | 14.621 | 184  | 146879   | 17.263 | ng/ul | 96       |
| 55) 4-Nitrophenol             | 14.721 | 109  | 185638   | 20.790 | ng/ul | 96       |
| 56) Dibenzofuran              | 14.892 | 168  | 1639445  | 21.376 | ng/ul | 99       |
| 57) 2,4-Dinitrotoluene        | 14.869 | 165  | 393961   | 21.221 | ng/ul | 96       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.116 | 232  | 326473   | 20.459 | ng/ul | 97       |
| 59) Diethylphthalate          | 15.316 | 149  | 1364743  | 21.175 | ng/ul | 99       |
| 61) Fluorene                  | 15.539 | 166  | 1393691  | 21.482 | ng/ul | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.533 | 204  | 682678   | 21.412 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.574 | 138  | 295118m  | 23.216 | ng/ul |          |
| 66) 4,6-Dinitro-2-methylph... | 15.627 | 198  | 230013   | 18.867 | ng/ul | 97       |
| 67) N-Nitrosodiphenylamine    | 15.751 | 169  | 1152309  | 21.454 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.427 | 248  | 400479   | 20.777 | ng/ul | 99       |
| 69) Hexachlorobenzene         | 16.533 | 284  | 490221   | 20.839 | ng/ul | 99       |
| 70) Atrazine                  | 16.704 | 200  | 382145   | 20.098 | ng/ul | 99       |
| 71) Pentachlorophenol         | 16.886 | 266  | 273592   | 19.005 | ng/ul | 97       |
| 72) Phenanthrene              | 17.274 | 178  | 2154499  | 21.267 | ng/ul | 100      |
| 74) Anthracene                | 17.362 | 178  | 2189265  | 21.478 | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.292 | 216  | 552072   | 21.027 | ng/uL | 99       |
| 76) Pentachlorobenzene        | 14.804 | 250  | 604342   | 21.154 | ng/uL | 99       |
| 77) Carbazole                 | 17.639 | 167  | 1936798  | 21.062 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 18.198 | 149  | 2268375  | 21.411 | ng/ul | 100      |
| 80) Fluoranthene              | 19.286 | 202  | 2411503  | 21.636 | ng/ul | 99       |
| 82) Pyrene                    | 19.651 | 202  | 2569918  | 21.887 | ng/ul | 100      |
| 83) Butylbenzylphthalate      | 20.545 | 149  | 985254   | 21.604 | ng/ul | 100      |
| 84) 3,3'-Dichlorobenzidine    | 21.327 | 252  | 841031   | 20.936 | ng/ul | 99       |
| 85) Benzo(a)anthracene        | 21.392 | 228  | 2409967  | 21.665 | ng/ul | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.315 | 149  | 1378766  | 21.682 | ng/ul | 100      |
| 87) Chrysene                  | 21.445 | 228  | 2270673  | 21.694 | ng/ul | 100      |
| 89) Di-n-octyl phthalate      | 22.221 | 149  | 2119373  | 20.809 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.039 | 252  | 2117494  | 21.744 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.086 | 252  | 2057887  | 21.583 | ng/ul | 99       |
| 93) Benzo(a)pyrene            | 23.650 | 252  | 1750847  | 21.497 | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.185 | 276  | 1846394  | 21.183 | ng/ul | 99       |
| 95) Dibenzo(a,h)anthracene    | 26.197 | 278  | 1663741  | 21.225 | ng/ul | 99       |
| 96) Benzo(g,h,i)perylene      | 26.932 | 276  | 1591535  | 21.310 | ng/ul | 99       |

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

**Instrument :**

BNM

**LabSampleId :**

SSTDCCC020

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/03/2022  
Supervised By :mohammad ahmed 11/04/2022

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110222\  
 Data File : BM037319.D  
 Acq On : 02 Nov 2022 16:53  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Nov 02 22:54:19 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 02 03:33:37 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2022  
 Supervised By :mohammad ahmed 11/04/2022

