

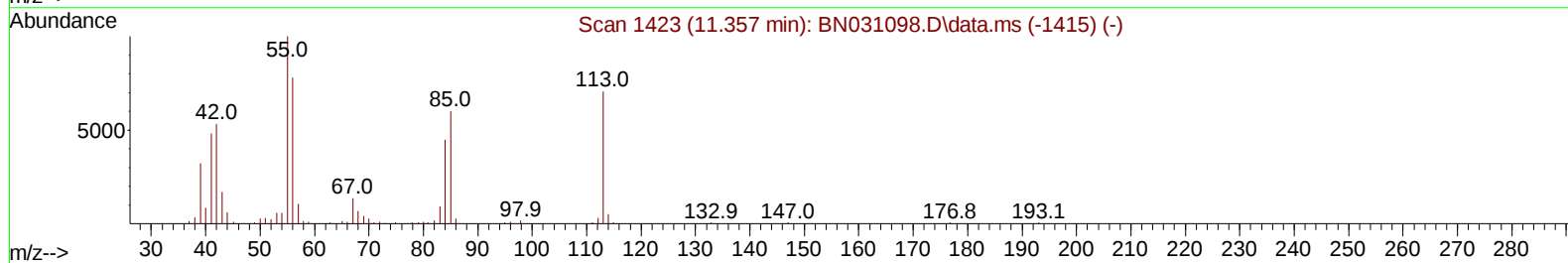
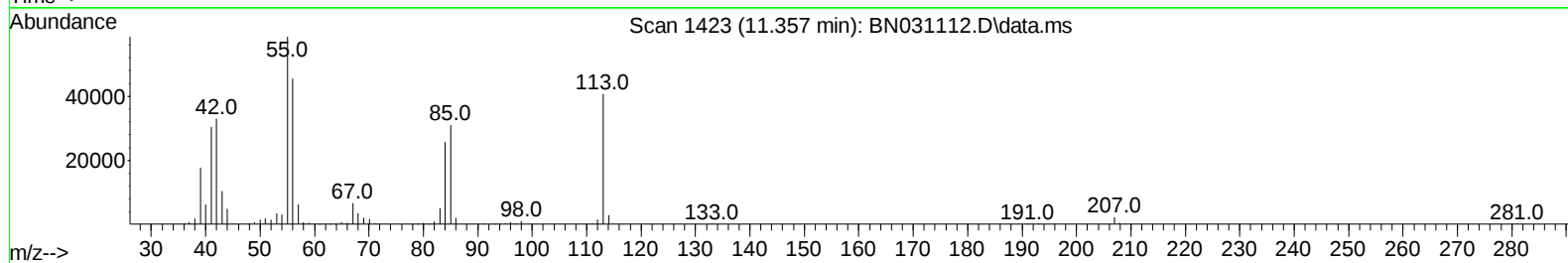
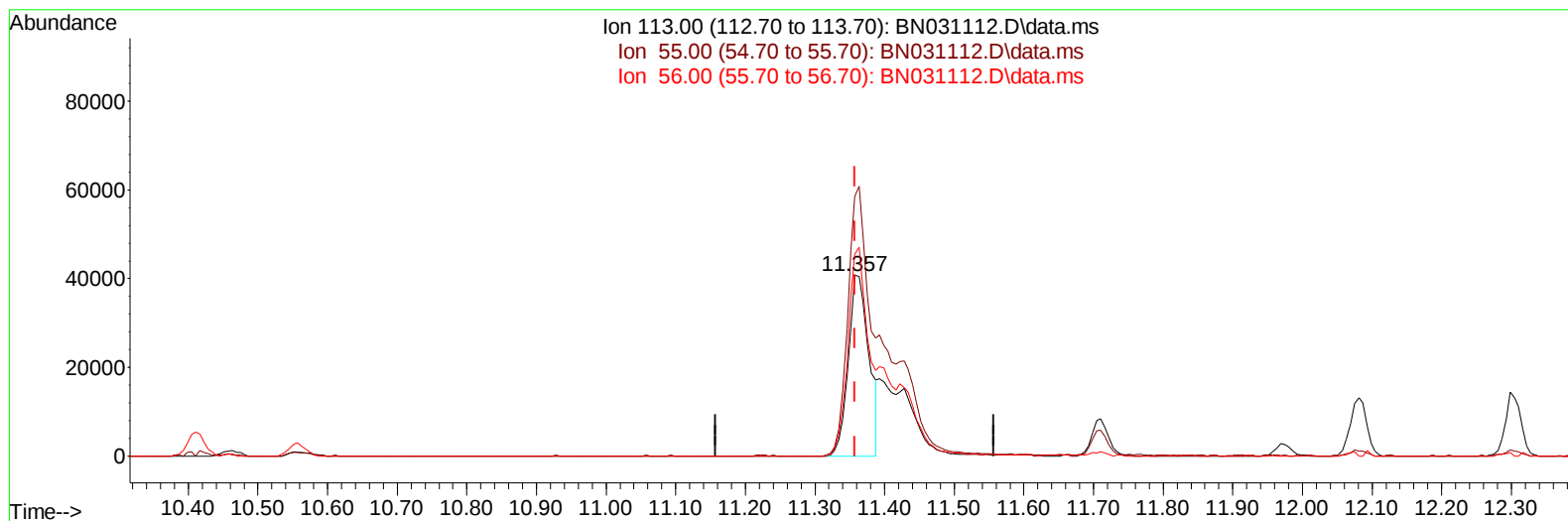
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041924\  
Data File : BN031112.D  
Acq On : 21 Apr 2024 17:06  
Operator : MA/JU  
Sample : PB160348BS  
Misc :  
ALS Vial : 63 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SLCS348

Manual IntegrationsAPPROVED

Quant Time: Apr 21 22:22:28 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Apr 19 22:56:25 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/22/2024  
Supervised By :mohammad ahmed 04/22/2024



TIC: BN031112.D\data.ms

(34) Caprolactam

11.357min (+ 0.000) 17.51 ng/ul

response 83737

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 140.50 | 143.88 |
| 56.00  | 110.00 | 111.61 |
| 0.00   | 0.00   | 0.00   |

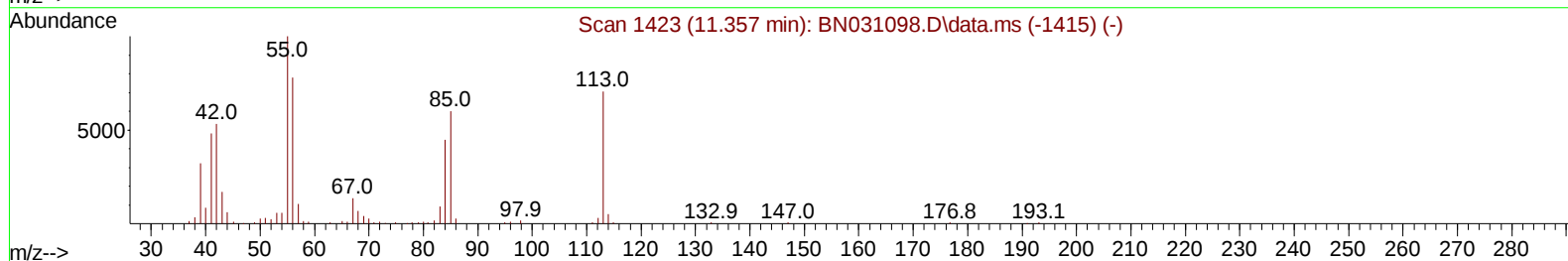
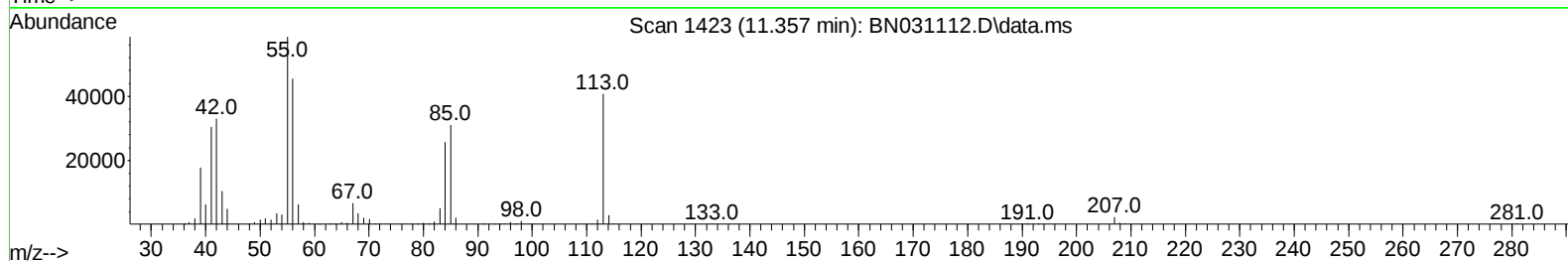
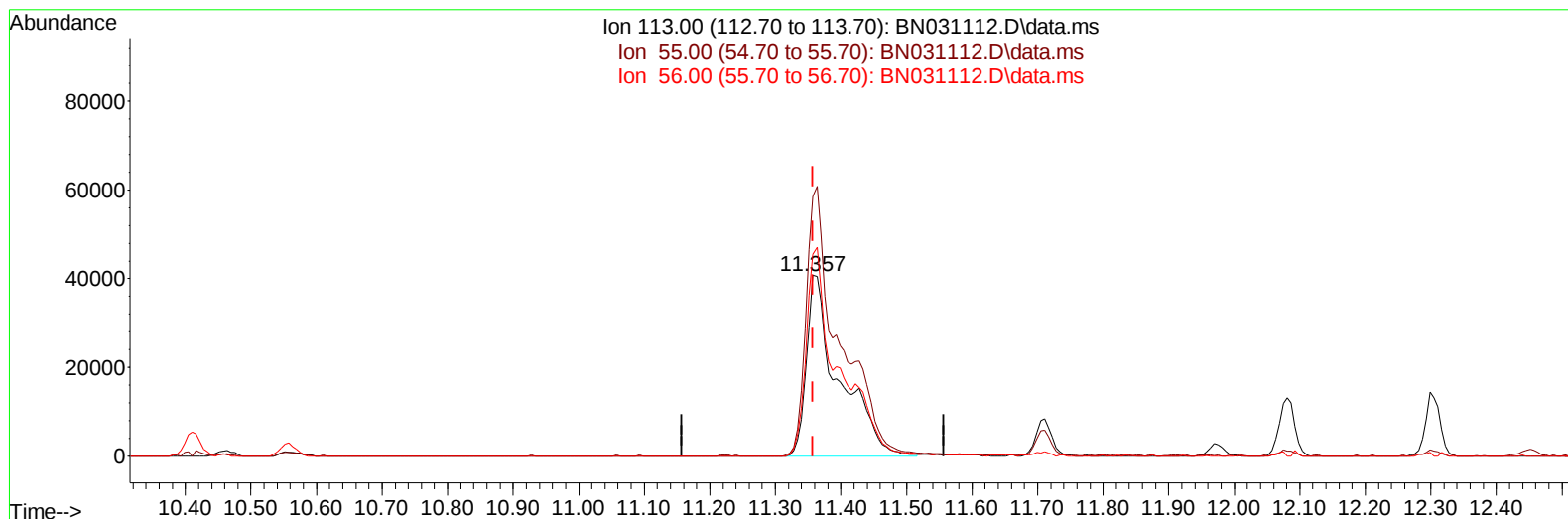
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041924\  
Data File : BN031112.D  
Acq On : 21 Apr 2024 17:06  
Operator : MA/JU  
Sample : PB160348BS  
Misc :  
ALS Vial : 63 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SLCS348

Manual IntegrationsAPPROVED

Quant Time: Apr 21 22:22:28 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Apr 19 22:56:25 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/22/2024  
Supervised By :mohammad ahmed 04/22/2024



TIC: BN031112.D\data.ms

(34) Caprolactam

11.357min (+ 0.000) 29.18 ng/ul m

response 139594

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 140.50 | 143.88 |
| 56.00  | 110.00 | 111.61 |
| 0.00   | 0.00   | 0.00   |

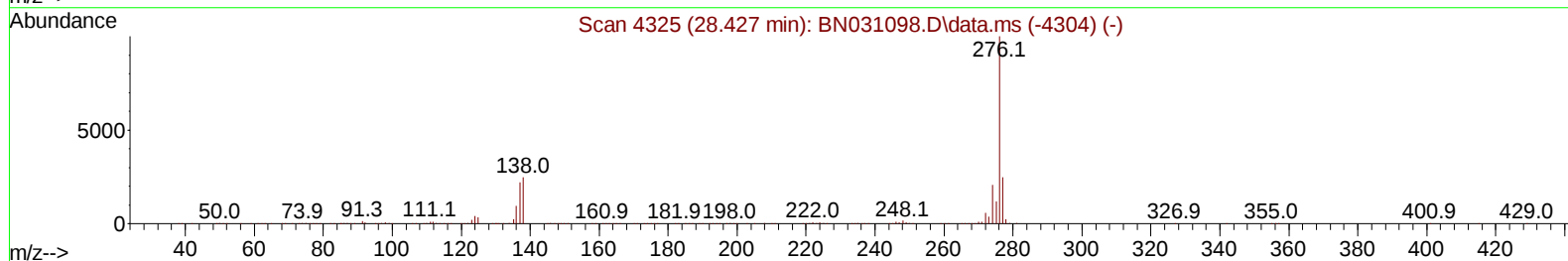
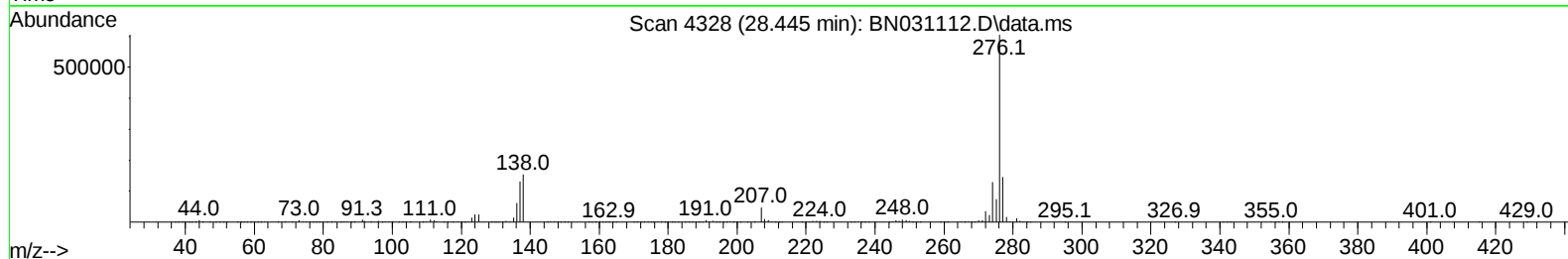
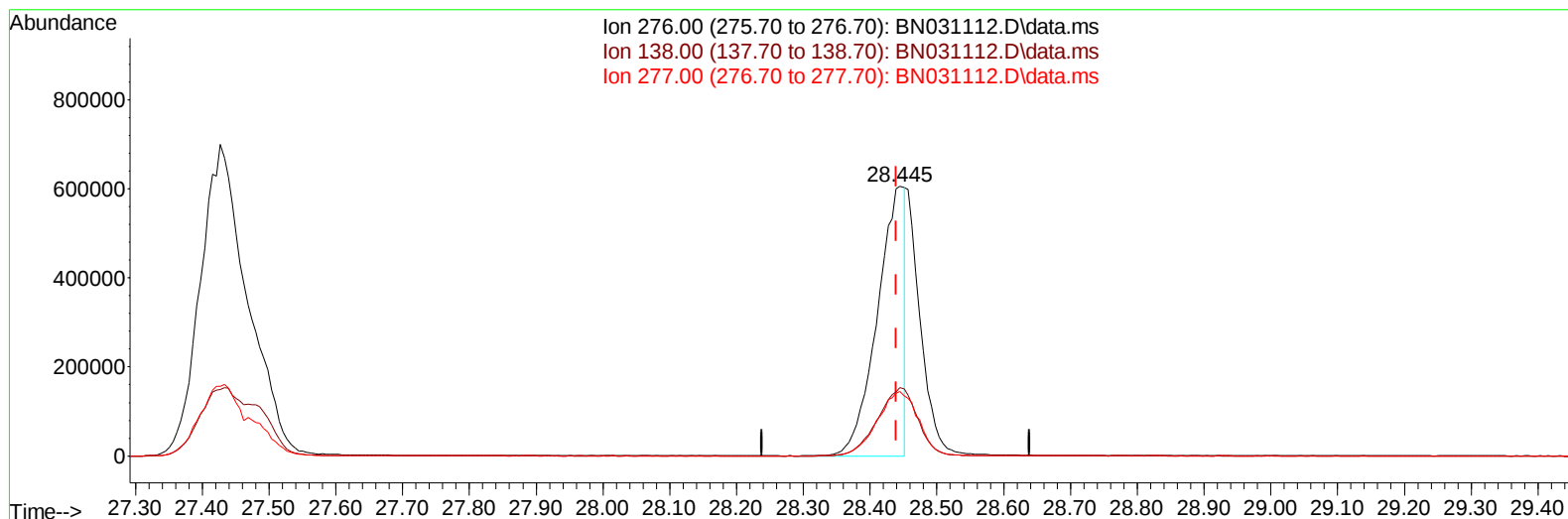
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041924\  
Data File : BN031112.D  
Acq On : 21 Apr 2024 17:06  
Operator : MA/JU  
Sample : PB160348BS  
Misc :  
ALS Vial : 63 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SLCS348

Manual IntegrationsAPPROVED

Quant Time: Apr 21 22:22:28 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Apr 19 22:56:25 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/22/2024  
Supervised By :mohammad ahmed 04/22/2024



TIC: BN031112.D\data.ms

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

28.445min (+ 0.006) 18.58 ng/u1

response 1713255

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 23.80  | 25.37  |
| 277.00 | 24.10  | 23.99  |
| 0.00   | 0.00   | 0.00   |

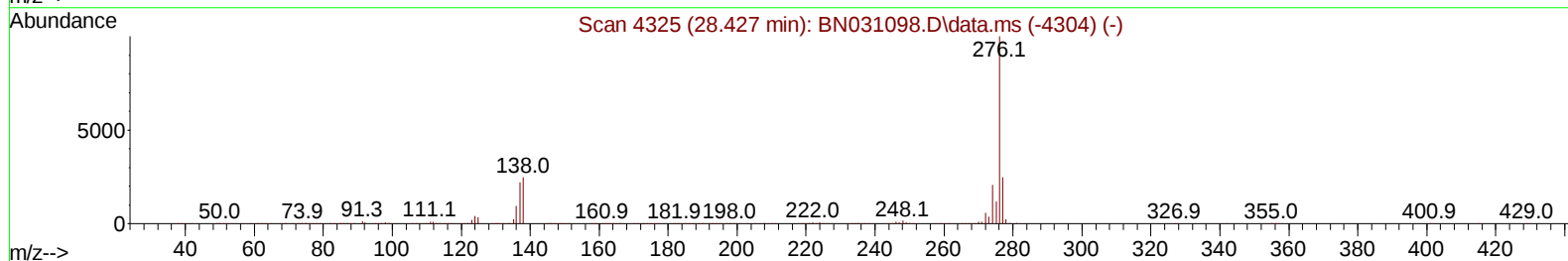
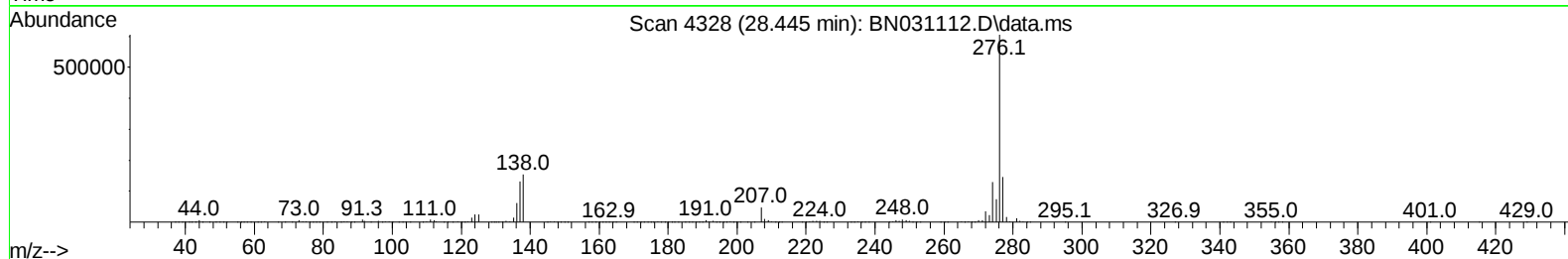
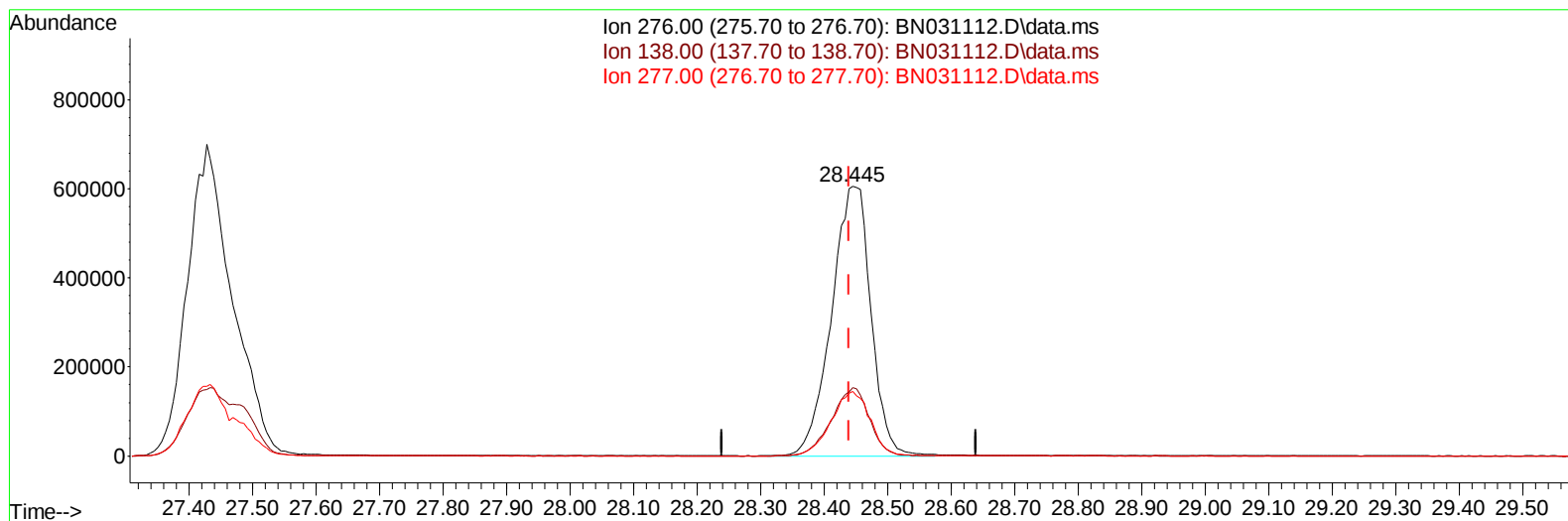
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041924\  
Data File : BN031112.D  
Acq On : 21 Apr 2024 17:06  
Operator : MA/JU  
Sample : PB160348BS  
Misc :  
ALS Vial : 63 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SLCS348

Manual IntegrationsAPPROVED

Quant Time: Apr 21 22:22:28 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Apr 19 22:56:25 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/22/2024  
Supervised By :mohammad ahmed 04/22/2024



TIC: BN031112.D\data.ms

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

28.445min (+ 0.006) 28.37 ng/ul m

response 2615548

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 23.80  | 25.37  |
| 277.00 | 24.10  | 23.99  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041924\  
 Data File : BN031112.D  
 Acq On : 21 Apr 2024 17:06  
 Operator : MA/JU  
 Sample : PB160348BS  
 Misc :  
 ALS Vial : 63 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SLCS348

Manual IntegrationsAPPROVED

Quant Time: Apr 21 22:33:05 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 19 22:56:25 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/22/2024  
 Supervised By :mohammad ahmed 04/22/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.628  | 152  | 180658   | 20.000 | ng/ul  | -0.01    |
| 20) Naphthalene-d8            | 10.410 | 136  | 825266   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.281 | 164  | 560277   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.034 | 188  | 1239689  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.269 | 240  | 1315842  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.174 | 264  | 1643567  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.146  | 96   | 26499    | 6.853  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.546  | 84   | 280891   | 25.223 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.805  | 99   | 466256   | 30.595 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.975  | 67   | 263959   | 29.770 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.164  | 132  | 364982   | 30.759 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.340  | 113  | 376290   | 29.360 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.787  | 128  | 187635   | 31.644 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.505  | 143  | 201512   | 32.694 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.034 | 165  | 380598   | 31.313 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.557 | 131  | 443230   | 23.335 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.698 | 166  | 1203670  | 30.873 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.969 | 160  | 1381098  | 31.260 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.493 | 143  | 259371   | 31.873 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.275 | 176  | 1048019  | 30.878 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.404 | 200  | 191122   | 30.558 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.128 | 188  | 1660327  | 30.900 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.433 | 212  | 2103127  | 32.526 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.974 | 264  | 2564364  | 32.902 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.175  | 88   | 38434    | 9.379  | ng/ul# | 91       |
| 5) Pyridine                   | 3.570  | 79   | 321039   | 28.569 | ng/ul  | 95       |
| 6) Benzaldehyde               | 6.781  | 77   | 258034   | 34.483 | ng/ul  | 98       |
| 8) Phenol                     | 6.834  | 94   | 470993   | 29.977 | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.069  | 93   | 353820   | 28.541 | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.199  | 128  | 367972   | 29.834 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.075  | 108  | 352416   | 28.944 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.158  | 45   | 466503   | 28.863 | ng/ul  | 99       |
| 16) Acetophenone              | 8.458  | 105  | 564591   | 28.432 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.440  | 70   | 278437   | 28.131 | ng/ul  | 99       |
| 18) 4-Methylphenol            | 8.405  | 108  | 400811   | 29.388 | ng/ul  | 97       |
| 19) Hexachloroethane          | 8.699  | 117  | 143898   | 28.116 | ng/ul  | 99       |
| 22) Nitrobenzene              | 8.834  | 77   | 417699   | 29.847 | ng/ul  | 98       |
| 23) Isophorone                | 9.352  | 82   | 812172   | 28.438 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.534  | 139  | 216514   | 31.750 | ng/ul  | 95       |
| 26) 2,4-Dimethylphenol        | 9.599  | 107  | 343982   | 24.734 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 9.840  | 93   | 501035   | 28.883 | ng/ul  | 98       |
| 29) 2,4-Dichlorophenol        | 10.063 | 162  | 378226   | 31.155 | ng/ul  | 99       |
| 30) Naphthalene               | 10.463 | 128  | 1233237  | 28.834 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.581 | 127  | 487245   | 26.401 | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 10.740 | 225  | 213990   | 28.914 | ng/ul  | 96       |
| 34) Caprolactam               | 11.357 | 113  | 139594m  | 29.184 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.710 | 107  | 428217   | 30.422 | ng/ul  | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041924\  
 Data File : BN031112.D  
 Acq On : 21 Apr 2024 17:06  
 Operator : MA/JU  
 Sample : PB160348BS  
 Misc :  
 ALS Vial : 63 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SLCS348

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/22/2024  
 Supervised By :mohammad ahmed 04/22/2024

Quant Time: Apr 21 22:33:05 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041924.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 19 22:56:25 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.081 | 142  | 878570   | 29.129 | ng/ul | 97       |
| 37) 1-Methylnaphthalene       | 12.304 | 142  | 897767   | 29.417 | ng/ul | 96       |
| 39) 1,2,4,5-Tetrachloroben... | 12.451 | 216  | 440008   | 30.013 | ng/ul | 97       |
| 40) Hexachlorocyclopentadiene | 12.428 | 237  | 171923   | 21.348 | ng/ul | 93       |
| 41) 2,4,6-Trichlorophenol     | 12.698 | 196  | 296348   | 30.368 | ng/ul | 88       |
| 42) 2,4,5-Trichlorophenol     | 12.769 | 196  | 337280   | 31.174 | ng/ul | 97       |
| 43) 1,1'-Biphenyl             | 13.110 | 154  | 1156671  | 30.017 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.146 | 162  | 905027   | 29.392 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.363 | 65   | 275389   | 32.397 | ng/ul | 99       |
| 47) Dimethylphthalate         | 13.746 | 163  | 1176736  | 29.256 | ng/ul | 99       |
| 48) 2,6-Dinitrotoluene        | 13.869 | 165  | 238369   | 31.571 | ng/ul | 98       |
| 50) Acenaphthylene            | 13.998 | 152  | 1485072  | 28.776 | ng/ul | 99       |
| 51) 3-Nitroaniline            | 14.198 | 138  | 277760   | 32.045 | ng/ul | 94       |
| 52) Acenaphthene              | 14.345 | 153  | 976082   | 28.531 | ng/ul | 97       |
| 53) 2,4-Dinitrophenol         | 14.404 | 184  | 116201   | 27.214 | ng/ul | 99       |
| 55) 4-Nitrophenol             | 14.510 | 109  | 204584   | 30.978 | ng/ul | 98       |
| 56) Dibenzofuran              | 14.681 | 168  | 1392514  | 29.439 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 14.657 | 165  | 360572   | 31.430 | ng/ul | 99       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.910 | 232  | 274379   | 29.344 | ng/ul | 99       |
| 59) Diethylphthalate          | 15.116 | 149  | 1196196  | 28.918 | ng/ul | 99       |
| 61) Fluorene                  | 15.334 | 166  | 1089955  | 28.000 | ng/ul | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.334 | 204  | 530233   | 28.597 | ng/ul | 98       |
| 63) 4-Nitroaniline            | 15.369 | 138  | 308946   | 33.059 | ng/ul | 95       |
| 66) 4,6-Dinitro-2-methylph... | 15.416 | 198  | 201602   | 30.019 | ng/ul | 100      |
| 67) N-Nitrosodiphenylamine    | 15.545 | 169  | 1015389  | 30.558 | ng/ul | 95       |
| 68) 4-Bromophenyl-phenylether | 16.228 | 248  | 344472   | 29.382 | ng/ul | 92       |
| 69) Hexachlorobenzene         | 16.328 | 284  | 391408   | 28.960 | ng/ul | 99       |
| 70) Atrazine                  | 16.504 | 200  | 260082   | 22.295 | ng/ul | 96       |
| 71) Pentachlorophenol         | 16.681 | 266  | 249748   | 28.441 | ng/ul | 98       |
| 72) Phenanthrene              | 17.075 | 178  | 1895326  | 28.905 | ng/ul | 98       |
| 74) Anthracene                | 17.163 | 178  | 1869417  | 28.172 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.069 | 216  | 446000   | 30.672 | ng/uL | 96       |
| 76) Pentachlorobenzene        | 14.598 | 250  | 438480   | 29.202 | ng/uL | 100      |
| 77) Carbazole                 | 17.439 | 167  | 1845955  | 30.455 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 18.004 | 149  | 2251345  | 30.472 | ng/ul | 99       |
| 80) Fluoranthene              | 19.098 | 202  | 2426603  | 31.709 | ng/ul | 99       |
| 82) Pyrene                    | 19.463 | 202  | 2535813  | 31.538 | ng/ul | 98       |
| 83) Butylbenzylphthalate      | 20.375 | 149  | 1134339  | 31.972 | ng/ul | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.186 | 252  | 837583   | 30.549 | ng/ul | 99       |
| 85) Benzo(a)anthracene        | 21.251 | 228  | 2748163  | 30.756 | ng/ul | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.180 | 149  | 1630140  | 30.601 | ng/ul | 100      |
| 87) Chrysene                  | 21.310 | 228  | 2639121  | 31.115 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.280 | 149  | 3124062  | 32.976 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.257 | 252  | 3025959  | 32.721 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.316 | 252  | 2960495  | 31.830 | ng/ul | 99       |
| 93) Benzo(a)pyrene            | 24.033 | 252  | 2569434  | 27.851 | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 27.427 | 276  | 3449199  | 30.129 | ng/ul | 99       |
| 95) Dibenzo(a,h)anthracene    | 27.480 | 278  | 2767170  | 29.449 | ng/ul | 98       |
| 96) Benzo(g,h,i)perylene      | 28.445 | 276  | 2615548m | 28.371 | ng/ul |          |

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

**Instrument :**

BNA\_N

**ClientSampleId :**

SLCS348

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 04/22/2024

Supervised By :mohammad ahmed 04/22/2024

