

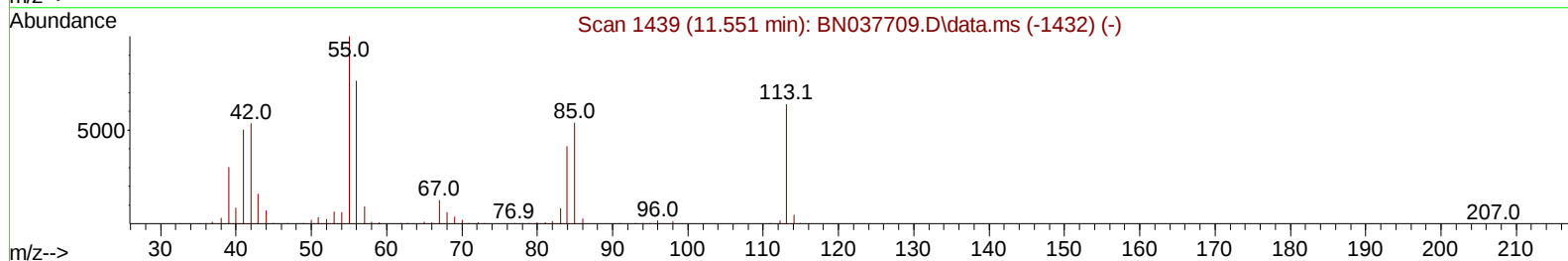
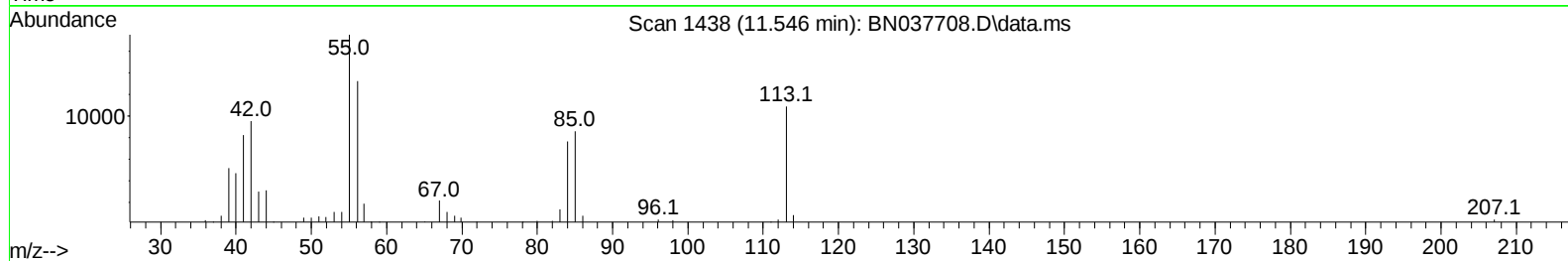
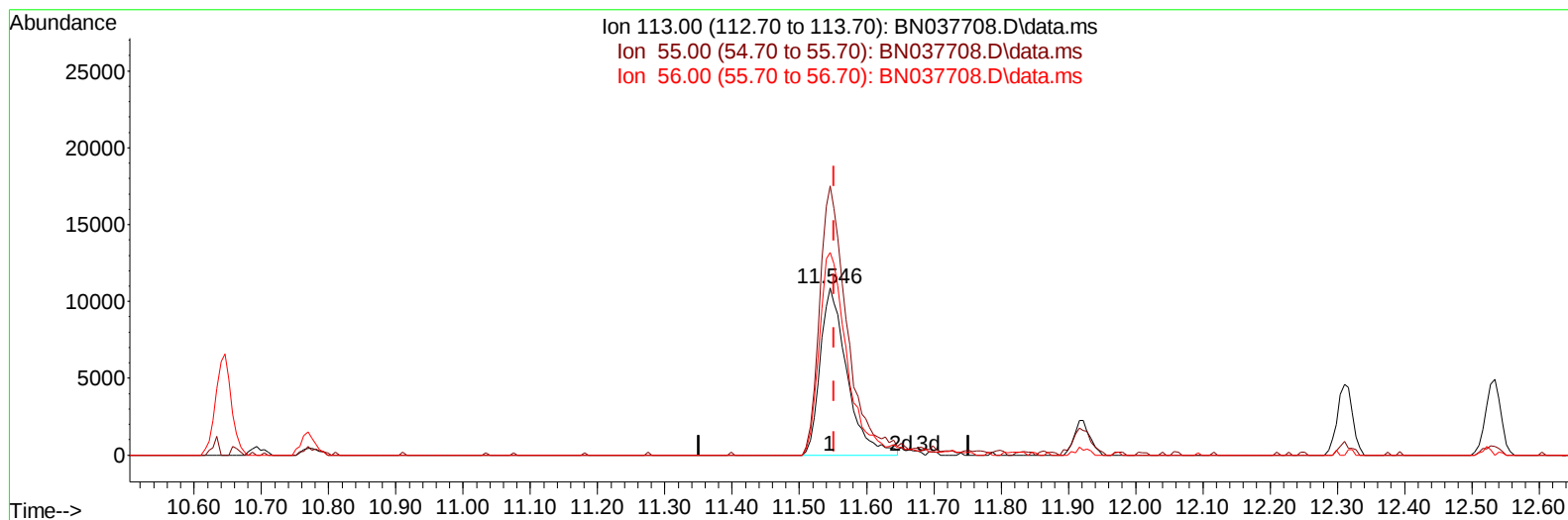
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091225\  
Data File : BN037708.D  
Acq On : 12 Sep 2025 12:08  
Operator : RC/JU  
Sample : SSTD01042  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Sep 12 15:26:04 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091225.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Sep 12 15:23:21 2025  
Response via : Initial Calibration

Reviewed By :Rahul Chavli 09/15/2025  
Supervised By :Jagrut Upadhyay 09/15/2025



TIC: BN037708.D\data.ms

(34) Caprolactam

11.546min (-0.006) 7.24 ng/u1

response 30087

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 158.30 | 161.00 |
| 56.00  | 120.50 | 121.27 |
| 0.00   | 0.00   | 0.00   |

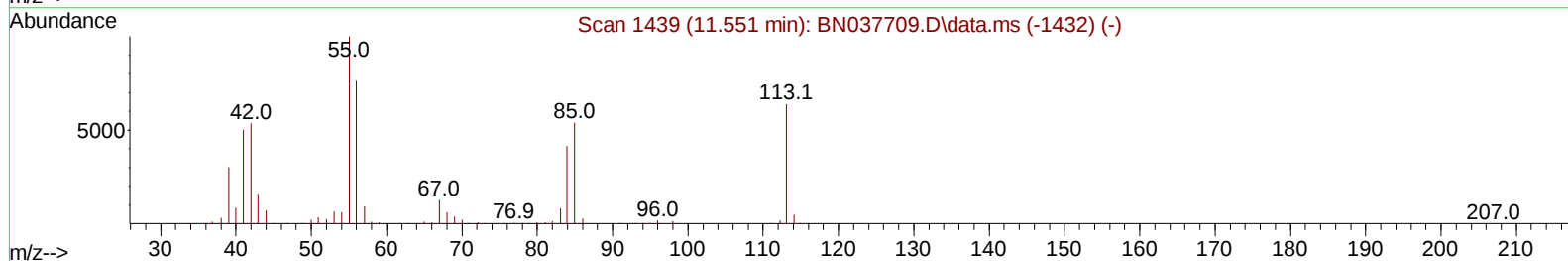
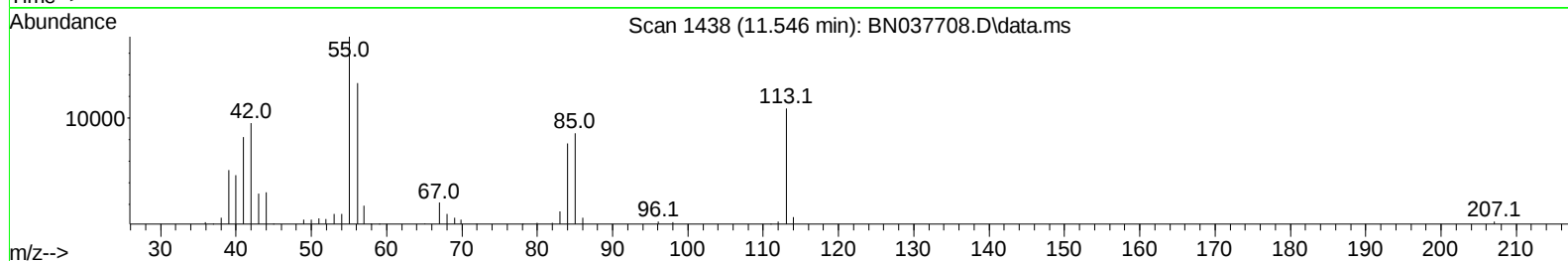
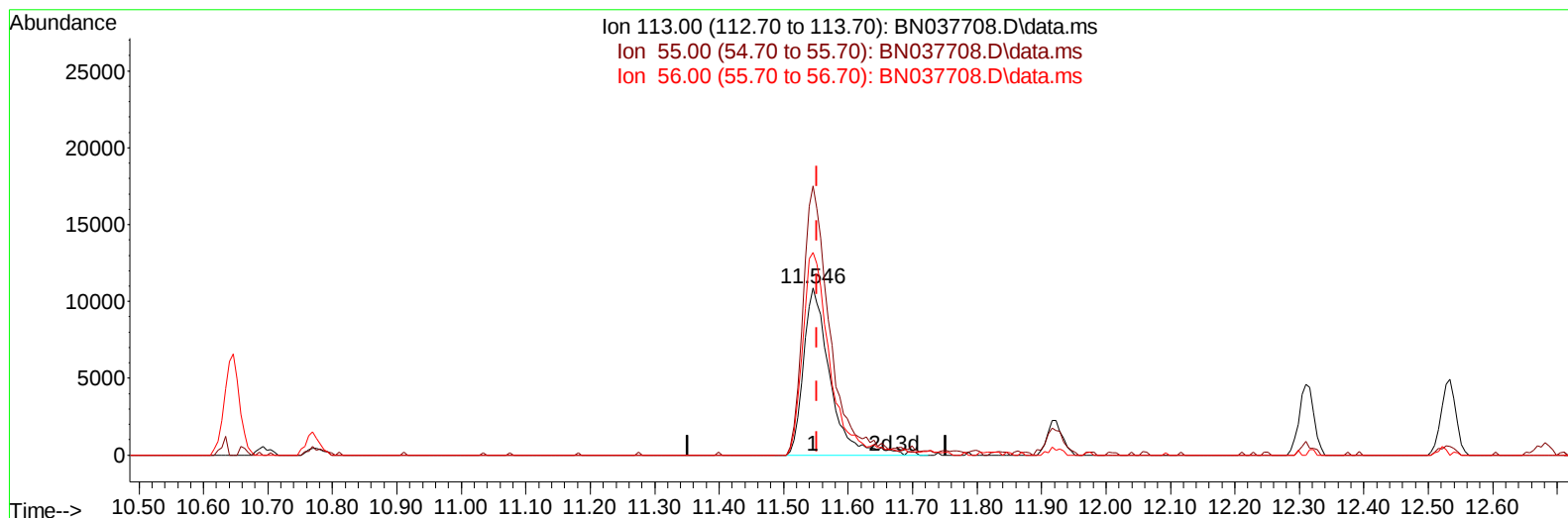
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091225\  
Data File : BN037708.D  
Acq On : 12 Sep 2025 12:08  
Operator : RC/JU  
Sample : SSTD01042  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Sep 12 15:26:04 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091225.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Sep 12 15:23:21 2025  
Response via : Initial Calibration

Reviewed By :Rahul Chavli 09/15/2025  
Supervised By :Jagrut Upadhyay 09/15/2025



TIC: BN037708.D\data.ms

**(34) Caprolactam**

11.546min (-0.006) 7.45 ng/ul m

response 30949

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 158.30 | 161.00 |
| 56.00  | 120.50 | 121.27 |
| 0.00   | 0.00   | 0.00   |

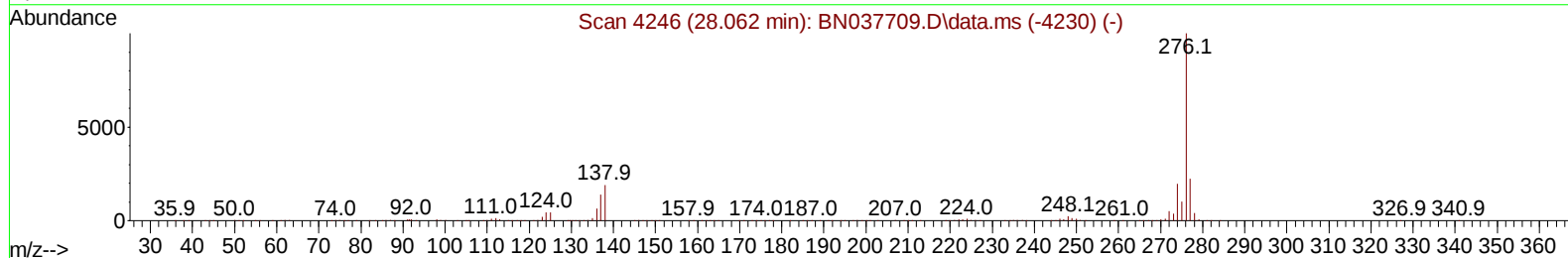
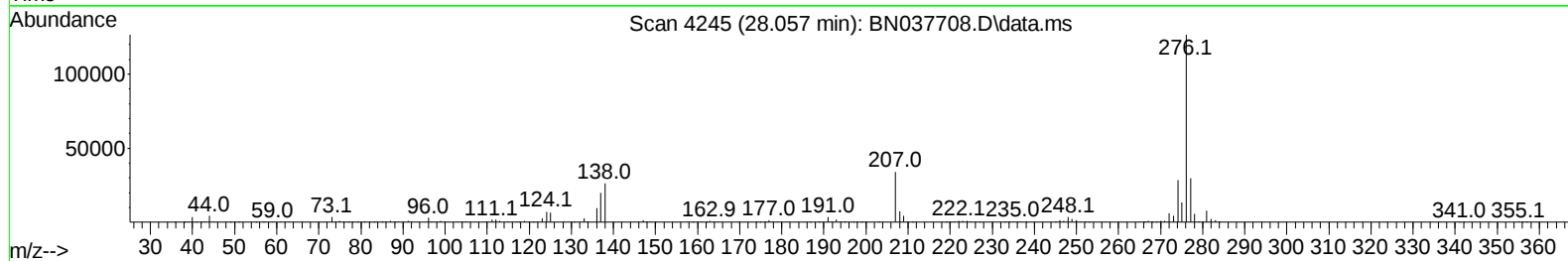
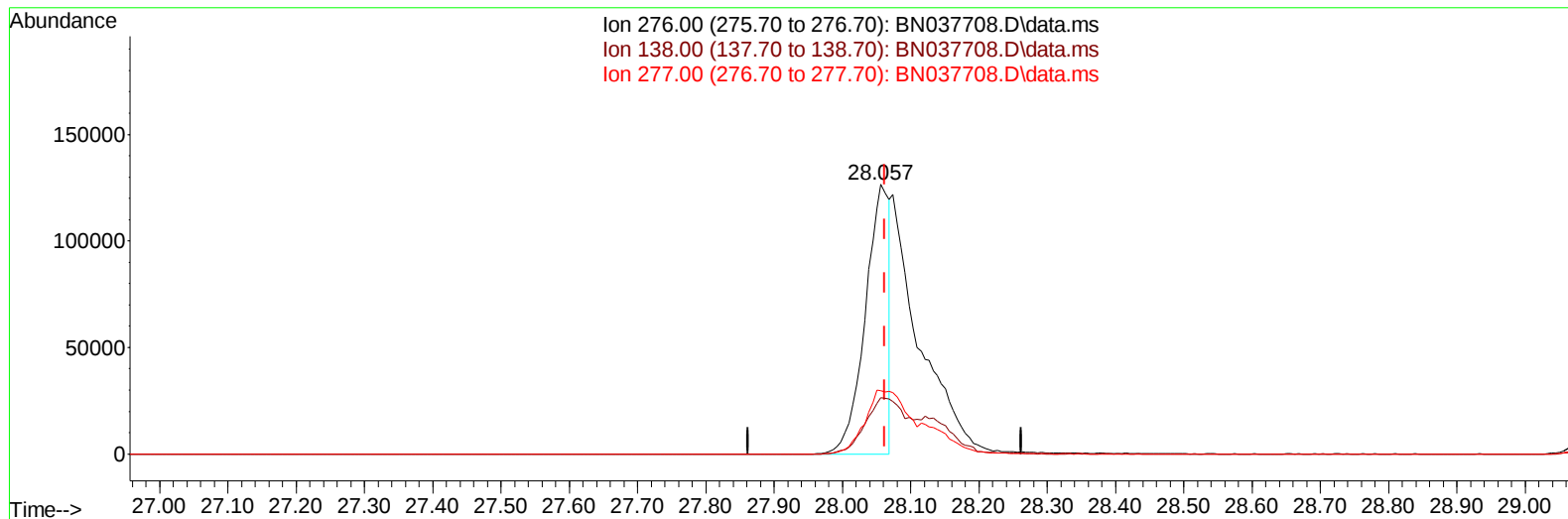
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091225\  
Data File : BN037708.D  
Acq On : 12 Sep 2025 12:08  
Operator : RC/JU  
Sample : SSTD01042  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Sep 12 15:26:04 2025  
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Quant Title : SVOA CALIBRATION  
QLast Update : Fri Sep 12 15:23:21 2025  
Response via : Initial Calibration

Reviewed By :Rahul Chavli 09/15/2025  
Supervised By :Jagrut Upadhyay 09/15/2025



TIC: BN037708.D\data.ms

**(94) Indeno(1,2,3-cd)pyrene**

**28.057min (-0.006) 4.38 ng/u1**

**response 307990**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 19.10  | 20.88  |
| 277.00 | 22.50  | 23.37  |
| 0.00   | 0.00   | 0.00   |

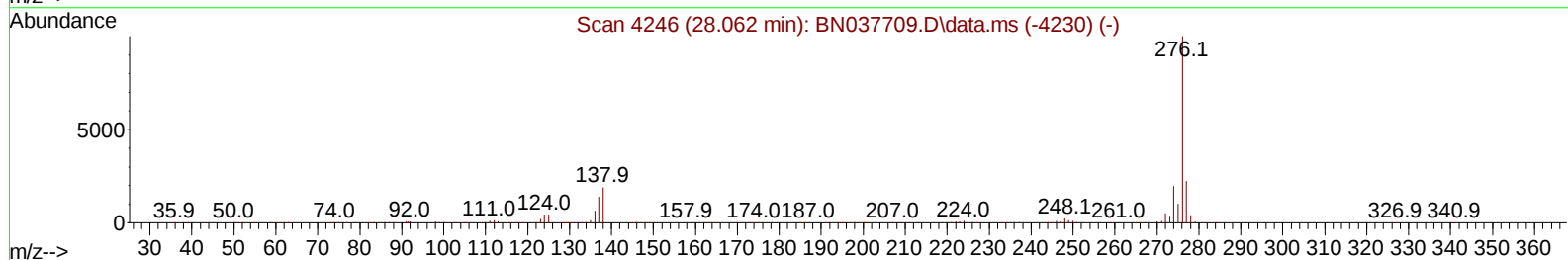
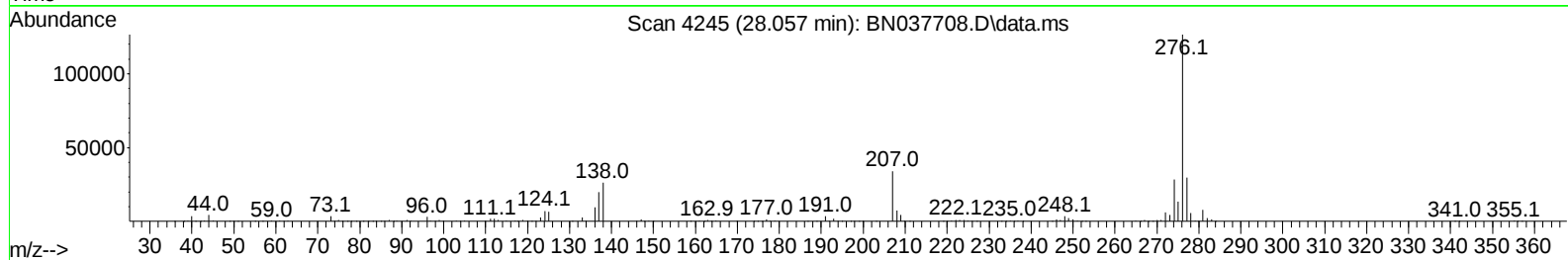
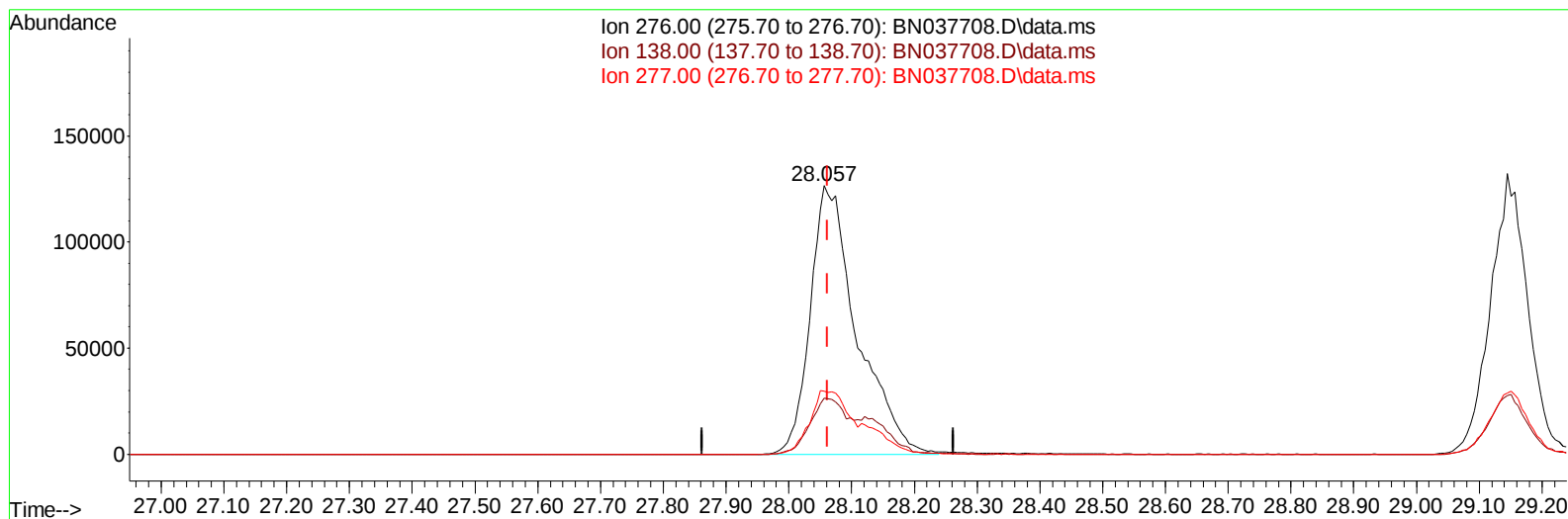
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091225\  
Data File : BN037708.D  
Acq On : 12 Sep 2025 12:08  
Operator : RC/JU  
Sample : SSTD01042  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Sep 12 15:26:04 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091225.MA.M  
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QLast Update : Fri Sep 12 15:23:21 2025  
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Reviewed By :Rahul Chavli 09/15/2025  
Supervised By :Jagrut Upadhyay 09/15/2025



TIC: BN037708.D\data.ms

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

28.057min (-0.006) 9.29 ng/ul m

response 653990

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 19.10  | 20.88  |
| 277.00 | 22.50  | 23.37  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091225\  
 Data File : BN037708.D  
 Acq On : 12 Sep 2025 12:08  
 Operator : RC/JU  
 Sample : SST001042  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST0010239

Manual IntegrationsAPPROVED

Quant Time: Sep 12 15:27:42 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091225.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 12 15:23:21 2025  
 Response via : Initial Calibration

Reviewed By :Rahul Chavli 09/15/2025  
 Supervised By :Jagrut Upadhyay 09/15/2025

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.834  | 152  | 239935   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.646 | 136  | 916337   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.493 | 164  | 533026   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.239 | 188  | 948495   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.492 | 240  | 905130   | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.580 | 264  | 962888   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 21761    | 4.187  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.675  | 84   | 146892   | 9.827  | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.987  | 99   | 165113   | 9.179  | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.158  | 67   | 108386   | 9.884  | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.364  | 132  | 139904   | 9.141  | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.534  | 113  | 129087   | 8.790  | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.993  | 128  | 64195    | 9.182  | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.716  | 143  | 45707    | 7.186  | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.257 | 165  | 121475   | 9.103  | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.769 | 131  | 186480   | 9.437  | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.898 | 166  | 342882   | 9.077  | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.187 | 160  | 429996   | 9.541  | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.669 | 143  | 47510    | 6.718  | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.487 | 176  | 309601   | 9.436  | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.592 | 200  | 24980    | 4.877  | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.339 | 188  | 452408   | 9.592  | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.628 | 212  | 469840   | 9.597  | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.369 | 264  | 436537   | 8.903  | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.299  | 88   | 23751    | 4.149  | ng/uL  | 94       |
| 5) Pyridine                   | 3.693  | 79   | 152289   | 9.917  | ng/ul  | 99       |
| 6) Benzaldehyde               | 6.964  | 77   | 106428   | 12.750 | ng/ul  | 98       |
| 8) Phenol                     | 7.011  | 94   | 176338   | 9.516  | ng/ul  | 100      |
| 10) Bis(2-Chloroethyl)ether   | 7.252  | 93   | 145231   | 9.683  | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.393  | 128  | 146313   | 9.320  | ng/ul  | 96       |
| 13) 2-Methylphenol            | 8.275  | 108  | 126455   | 9.033  | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.369  | 45   | 200689   | 9.295  | ng/ul  | 99       |
| 16) Acetophenone              | 8.652  | 105  | 220410   | 9.576  | ng/ul  | 100      |
| 17) N-Nitroso-di-n-propyla... | 8.640  | 70   | 99512    | 9.306  | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.599  | 108  | 139312   | 9.242  | ng/ul  | 99       |
| 19) Hexachloroethane          | 8.928  | 117  | 63429    | 9.536  | ng/ul  | 98       |
| 22) Nitrobenzene              | 9.034  | 77   | 152714   | 9.556  | ng/ul  | 100      |
| 23) Isophorone                | 9.563  | 82   | 266518   | 9.216  | ng/ul  | 98       |
| 25) 2-Nitrophenol             | 9.752  | 139  | 56660    | 7.780  | ng/ul  | 99       |
| 26) 2,4-Dimethylphenol        | 9.810  | 107  | 146331   | 9.459  | ng/ul  | 97       |
| 27) Bis(2-Chloroethoxy)met... | 10.052 | 93   | 181210   | 9.628  | ng/ul  | 98       |
| 29) 2,4-Dichlorophenol        | 10.281 | 162  | 124912   | 9.267  | ng/ul  | 98       |
| 30) Naphthalene               | 10.693 | 128  | 488651   | 9.890  | ng/ul  | 98       |
| 32) 4-Chloroaniline           | 10.793 | 127  | 188563   | 9.519  | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 10.993 | 225  | 91048    | 9.741  | ng/ul  | 99       |
| 34) Caprolactam               | 11.546 | 113  | 30949m   | 7.446  | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.922 | 107  | 118223   | 8.713  | ng/ul  | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091225\  
 Data File : BN037708.D  
 Acq On : 12 Sep 2025 12:08  
 Operator : RC/JU  
 Sample : SST001042  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST0010239

Manual IntegrationsAPPROVED

Quant Time: Sep 12 15:27:42 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091225.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 12 15:23:21 2025  
 Response via : Initial Calibration

Reviewed By :Rahul Chavli 09/15/2025  
 Supervised By :Jagrut Upadhyay 09/15/2025

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.310 | 142  | 322882   | 9.636  | ng/ul | 98       |
| 37) 1-Methylnaphthalene       | 12.534 | 142  | 315964   | 9.632  | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.681 | 216  | 159376   | 9.772  | ng/ul | 98       |
| 40) Hexachlorocyclopentadiene | 12.669 | 237  | 92018    | 8.818  | ng/ul | 98       |
| 41) 2,4,6-Trichlorophenol     | 12.916 | 196  | 80290    | 8.440  | ng/ul | 99       |
| 42) 2,4,5-Trichlorophenol     | 12.987 | 196  | 95468    | 8.810  | ng/ul | 99       |
| 43) 1,1'-Biphenyl             | 13.328 | 154  | 401236   | 9.806  | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.363 | 162  | 316888   | 9.970  | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.557 | 65   | 63029    | 8.402  | ng/ul | 95       |
| 47) Dimethylphthalate         | 13.951 | 163  | 344429   | 9.231  | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 14.057 | 165  | 55909    | 7.945  | ng/ul | 95       |
| 50) Acenaphthylene            | 14.216 | 152  | 504847   | 9.788  | ng/ul | 99       |
| 51) 3-Nitroaniline            | 14.381 | 138  | 60292    | 8.628  | ng/ul | 99       |
| 52) Acenaphthene              | 14.557 | 153  | 326965   | 9.777  | ng/ul | 95       |
| 53) 2,4-Dinitrophenol         | 14.587 | 184  | 14995    | 4.223  | ng/ul | 92       |
| 55) 4-Nitrophenol             | 14.681 | 109  | 41032    | 8.046  | ng/ul | 99       |
| 56) Dibenzofuran              | 14.893 | 168  | 452663   | 9.846  | ng/ul | 96       |
| 57) 2,4-Dinitrotoluene        | 14.845 | 165  | 79463    | 7.804  | ng/ul | 98       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.116 | 232  | 63927    | 7.745  | ng/ul | 99       |
| 59) Diethylphthalate          | 15.322 | 149  | 322850   | 8.701  | ng/ul | 99       |
| 61) Fluorene                  | 15.545 | 166  | 366585   | 9.782  | ng/ul | 95       |
| 62) 4-Chlorophenyl-phenyle... | 15.540 | 204  | 179570   | 9.799  | ng/ul | 98       |
| 63) 4-Nitroaniline            | 15.545 | 138  | 60727    | 9.269  | ng/ul | 95       |
| 66) 4,6-Dinitro-2-methylph... | 15.610 | 198  | 30071    | 5.433  | ng/ul | 96       |
| 67) N-Nitrosodiphenylamine    | 15.751 | 169  | 292992   | 9.718  | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.434 | 248  | 102949   | 9.596  | ng/ul | 93       |
| 69) Hexachlorobenzene         | 16.551 | 284  | 127555   | 10.011 | ng/ul | 97       |
| 70) Atrazine                  | 16.698 | 200  | 87043    | 8.201  | ng/ul | 97       |
| 71) Pentachlorophenol         | 16.887 | 266  | 49579    | 6.417  | ng/ul | 93       |
| 72) Phenanthrene              | 17.281 | 178  | 535956   | 9.839  | ng/ul | 98       |
| 74) Anthracene                | 17.375 | 178  | 538496   | 9.739  | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.287 | 216  | 159687   | 10.003 | ng/uL | 97       |
| 76) Pentachlorobenzene        | 14.816 | 250  | 156763   | 10.168 | ng/uL | 96       |
| 77) Carbazole                 | 17.639 | 167  | 462589   | 9.522  | ng/ul | 98       |
| 78) Di-n-butylphthalate       | 18.222 | 149  | 453361   | 7.917  | ng/ul | 99       |
| 80) Fluoranthene              | 19.298 | 202  | 543410   | 9.390  | ng/ul | 99       |
| 82) Pyrene                    | 19.657 | 202  | 595941   | 9.852  | ng/ul | 99       |
| 83) Butylbenzylphthalate      | 20.569 | 149  | 131929   | 6.471  | ng/ul | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.392 | 252  | 141807   | 8.048  | ng/ul | 97       |
| 85) Benzo(a)anthracene        | 21.475 | 228  | 567531   | 9.438  | ng/ul | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.416 | 149  | 202276   | 6.079  | ng/ul | 97       |
| 87) Chrysene                  | 21.533 | 228  | 540465   | 9.501  | ng/ul | 97       |
| 89) Di-n-octyl phthalate      | 22.592 | 149  | 308197   | 5.402  | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.604 | 252  | 536353   | 9.376  | ng/ul | 98       |
| 91) Benzo(k)fluoranthene      | 23.669 | 252  | 538266   | 9.098  | ng/ul | 100      |
| 93) Benzo(a)pyrene            | 24.427 | 252  | 519045   | 9.501  | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.057 | 276  | 653990m  | 9.292  | ng/ul |          |
| 95) Dibenzo(a,h)anthracene    | 28.139 | 278  | 539554   | 9.241  | ng/ul | 98       |
| 96) Benzo(g,h,i)perylene      | 29.145 | 276  | 548196   | 9.666  | ng/ul | 98       |

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

**Instrument :**

BNA\_N

**ClientSampleId :**

SSTD010239

**Manual IntegrationsAPPROVED**

Reviewed By :Rahul Chavli 09/15/2025

Supervised By :Jagrut Upadhyay 09/15/2025

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091225\  
 Data File : BN037708.D  
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 Operator : RC/JU  
 Sample : SSTD01042  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD010239

Manual IntegrationsAPPROVED

Quant Time: Sep 12 15:27:42 2025  
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