

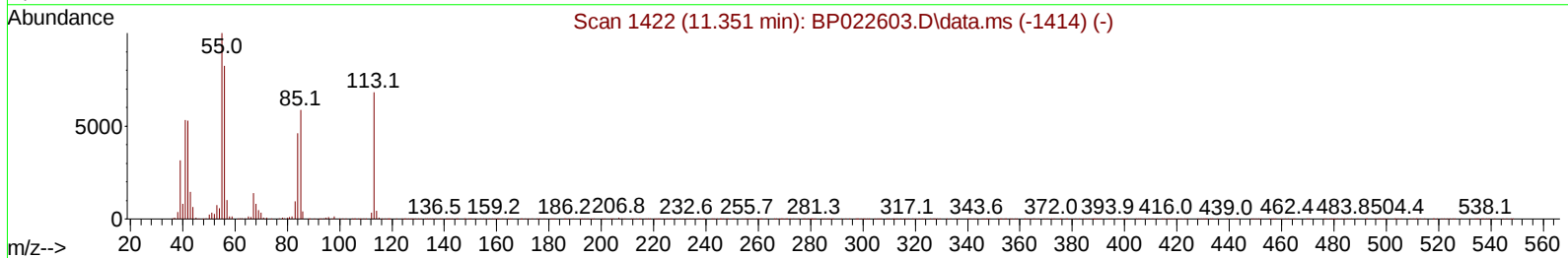
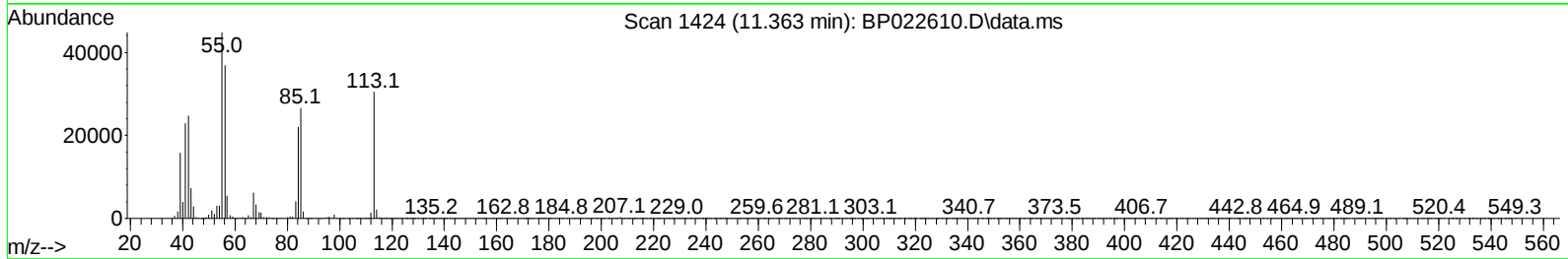
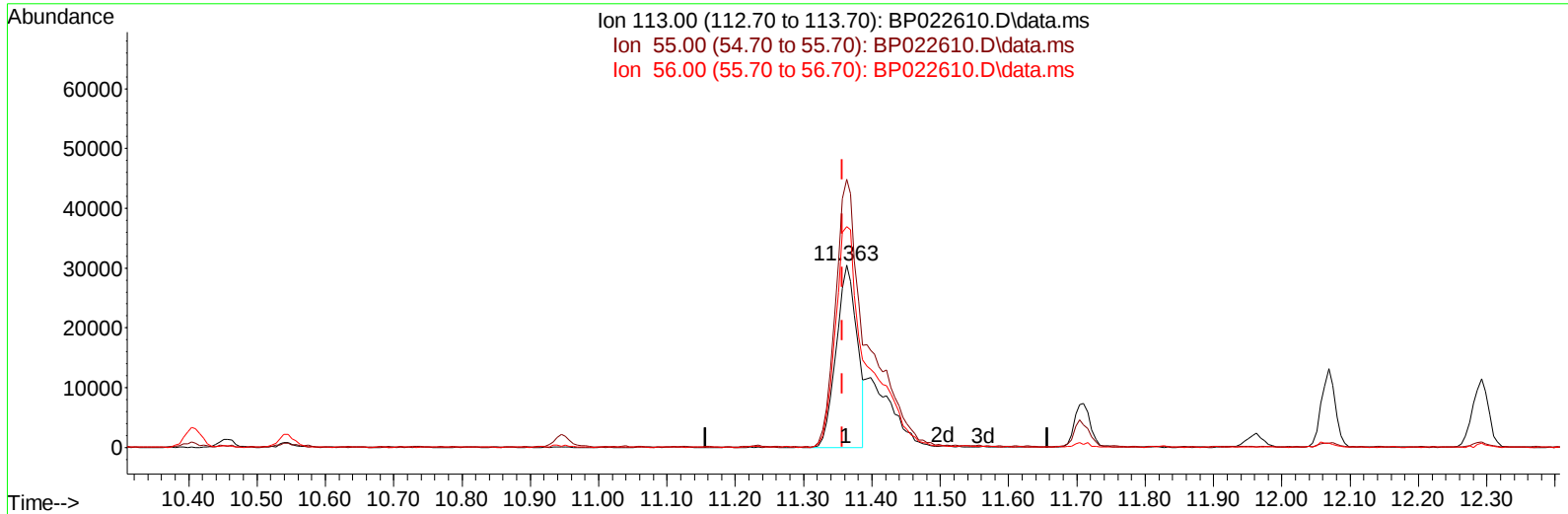
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102524\  
Data File : BP022610.D  
Acq On : 25 Oct 2024 13:55  
Operator : RC/JU  
Sample : P4435-03MSD  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BHK68MSD

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/26/2024  
Supervised By :mohammad ahmed 10/29/2024

Quant Time: Oct 25 14:50:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Oct 22 05:59:59 2024  
Response via : Initial Calibration



TIC: BP022610.D\data.ms

## (34) Caprolactam

11.363min (+ 0.006) 23.12 ng/ul

response 64725

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 162.70 | 146.83 |
| 56.00  | 130.00 | 120.86 |
| 0.00   | 0.00   | 0.00   |

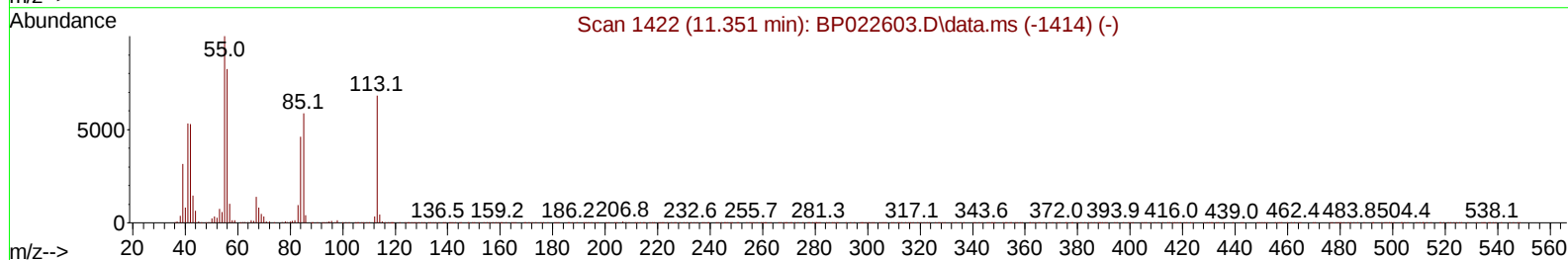
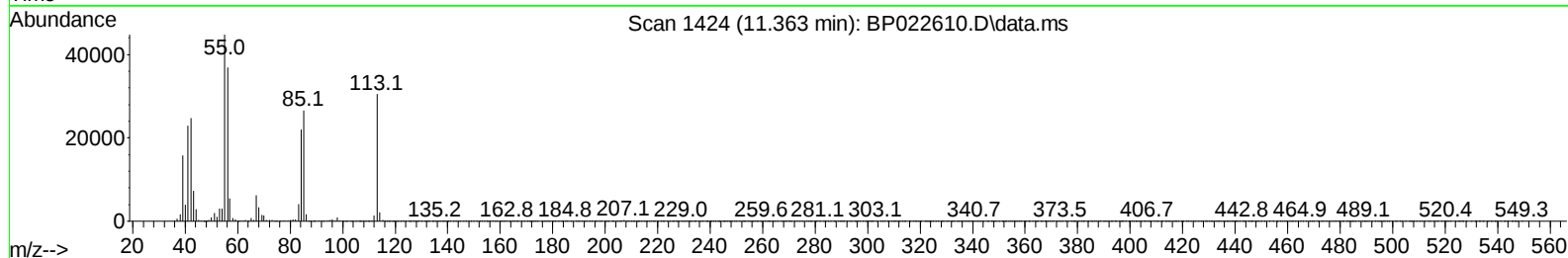
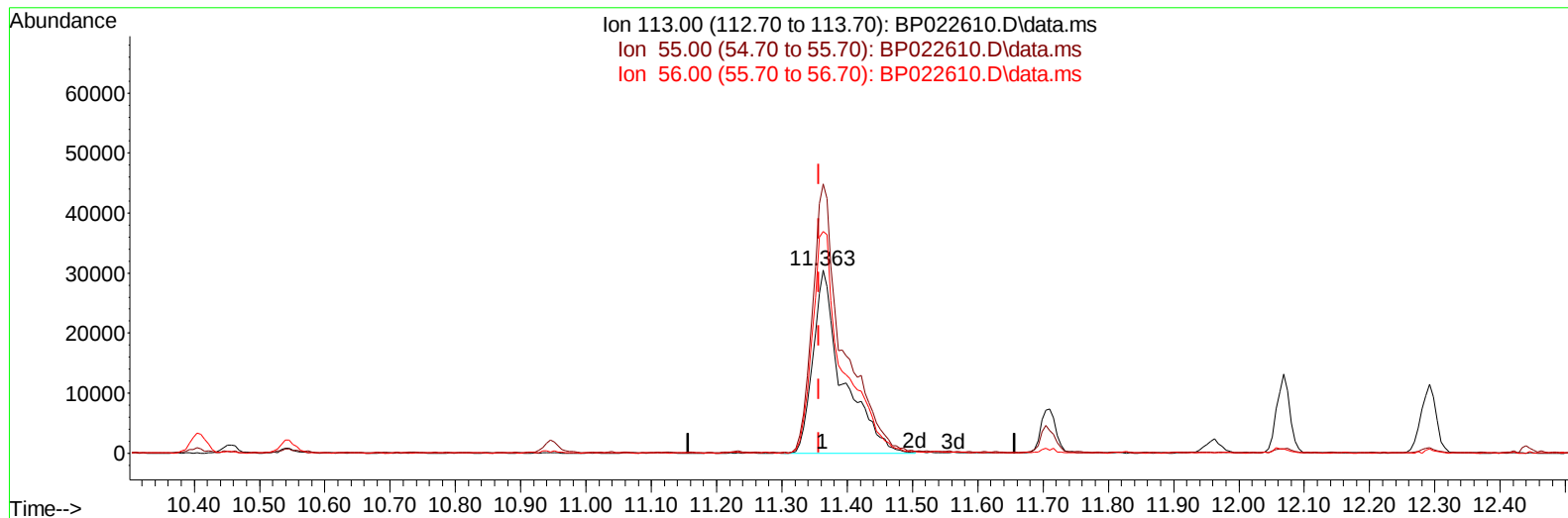
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102524\  
Data File : BP022610.D  
Acq On : 25 Oct 2024 13:55  
Operator : RC/JU  
Sample : P4435-03MSD  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BHK68MSD

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/26/2024  
Supervised By :mohammad ahmed 10/29/2024

Quant Time: Oct 25 14:50:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Oct 22 05:59:59 2024  
Response via : Initial Calibration



TIC: BP022610.D\data.ms

## (34) Caprolactam

11.363min (+ 0.006) 34.52 ng/ul m

response 96645

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 162.70 | 146.83 |
| 56.00  | 130.00 | 120.86 |
| 0.00   | 0.00   | 0.00   |

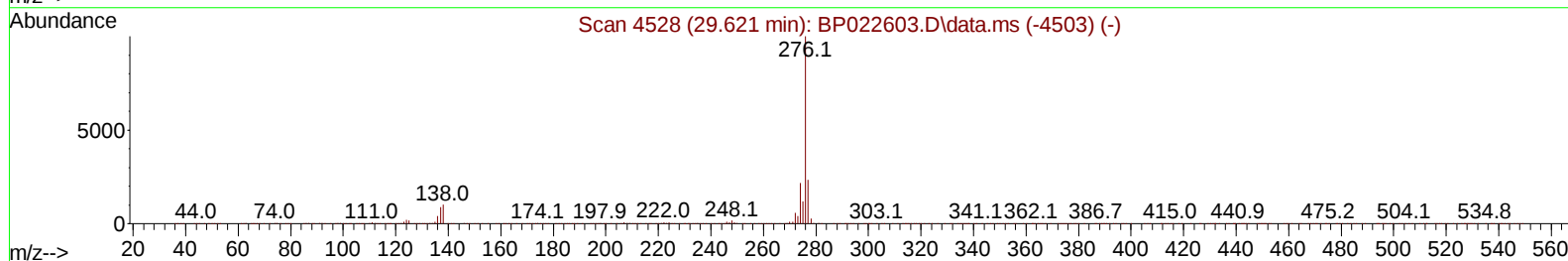
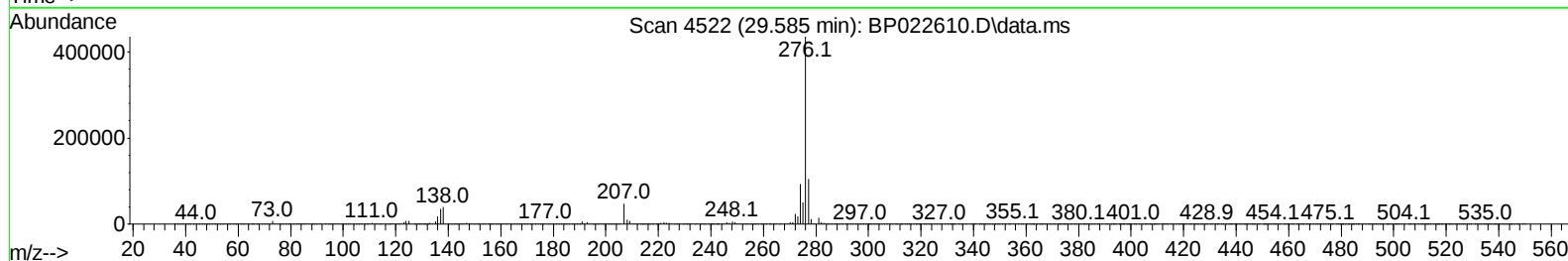
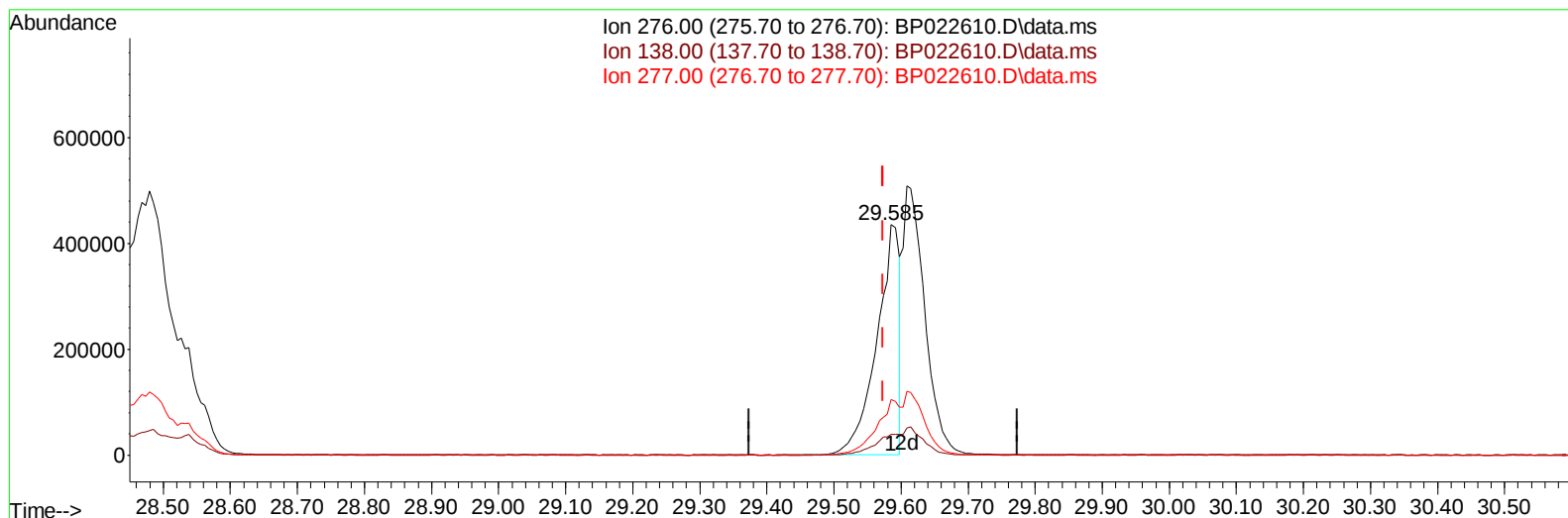
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102524\  
Data File : BP022610.D  
Acq On : 25 Oct 2024 13:55  
Operator : RC/JU  
Sample : P4435-03MSD  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BHK68MSD

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/26/2024  
Supervised By :mohammad ahmed 10/29/2024

Quant Time: Oct 25 14:50:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Oct 22 05:59:59 2024  
Response via : Initial Calibration



TIC: BP022610.D\data.ms

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

29.585min (+ 0.012) 15.91 ng/u1

response 1019525

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 10.60  | 9.01   |
| 277.00 | 23.70  | 24.03  |
| 0.00   | 0.00   | 0.00   |

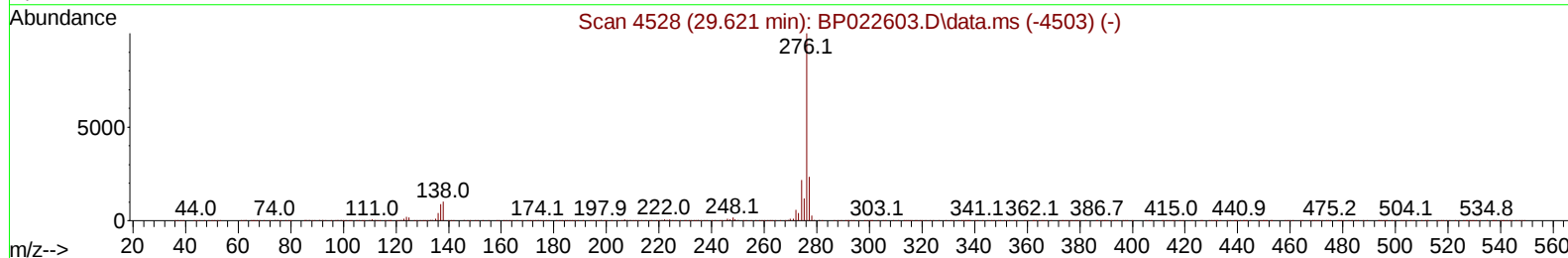
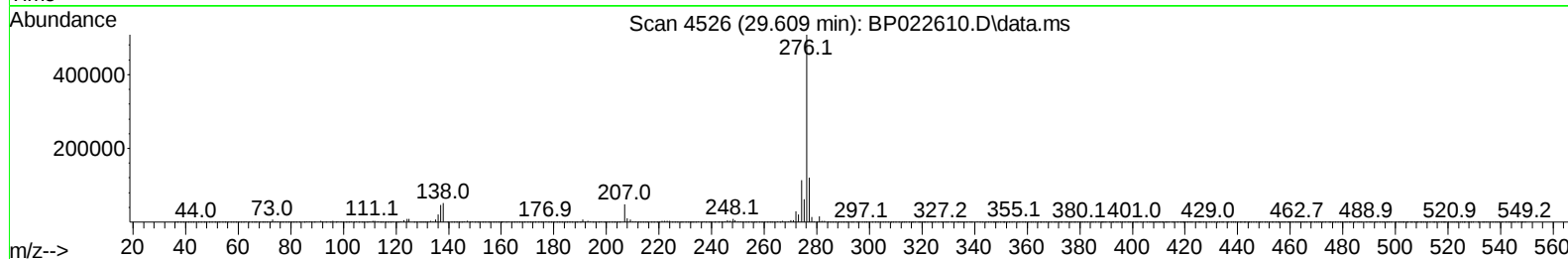
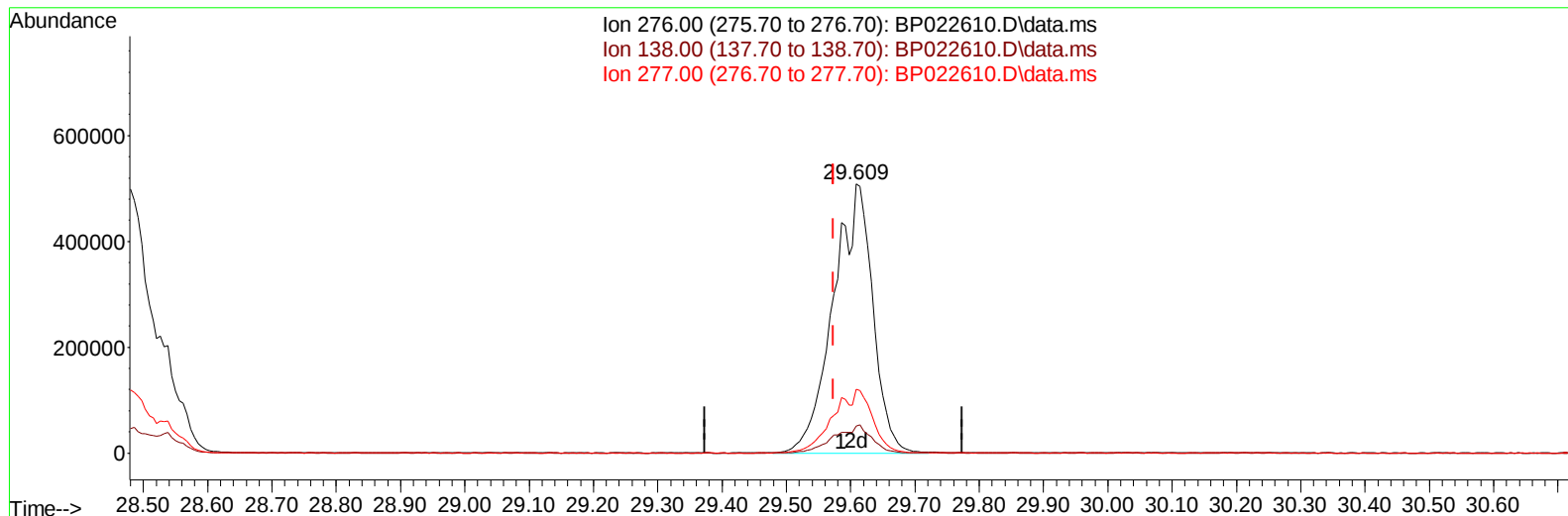
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102524\  
Data File : BP022610.D  
Acq On : 25 Oct 2024 13:55  
Operator : RC/JU  
Sample : P4435-03MSD  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BHK68MSD

Manual IntegrationsAPPROVED

Quant Time: Oct 25 14:50:32 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Oct 22 05:59:59 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/26/2024  
Supervised By :mohammad ahmed 10/29/2024



TIC: BP022610.D\data.ms

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

29.609min (+ 0.035) 33.90 ng/ul m

response 2172386

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 10.60  | 10.16  |
| 277.00 | 23.70  | 23.83  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102524\  
 Data File : BP022610.D  
 Acq On : 25 Oct 2024 13:55  
 Operator : RC/JU  
 Sample : P4435-03MSD  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHK68MSD

Manual IntegrationsAPPROVED

Quant Time: Oct 25 14:52:28 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 22 05:59:59 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/26/2024  
 Supervised By :mohammad ahmed 10/29/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.651  | 152  | 117045   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.410 | 136  | 466104   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.269 | 164  | 381007   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.063 | 188  | 932931   | 20.000 | ng/ul  | -0.02    |
| 79) Chrysene-d12              | 21.498 | 240  | 937089   | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.768 | 264  | 1065811  | 20.000 | ng/ul  | 0.02     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.187  | 96   | 12483    | 4.144  | ng/uL  | -0.01    |
| 4) Pyridine-d5                | 3.593  | 84   | 170771   | 20.653 | ng/ul  | -0.02    |
| 7) Phenol-d5                  | 6.845  | 99   | 279446   | 28.363 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.992  | 67   | 148885   | 25.711 | ng/ul  | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.192  | 132  | 200431   | 28.670 | ng/ul  | -0.01    |
| 15) 4-Methylphenol-d8         | 8.357  | 113  | 222832   | 28.340 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.792  | 128  | 101728   | 28.385 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.504  | 143  | 126286   | 30.436 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.039 | 165  | 267391   | 30.321 | ng/ul  | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.545 | 131  | 267595   | 25.680 | ng/ul  | -0.01    |
| 46) Dimethylphthalate-d6      | 13.680 | 166  | 837391   | 27.662 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.957 | 160  | 900251   | 28.000 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.510 | 143  | 117815   | 23.199 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.274 | 176  | 739485   | 28.164 | ng/ul  | -0.02    |
| 65) 4,6-Dinitro-2-methylph... | 15.427 | 200  | 141951   | 23.135 | ng/ul  | 0.01     |
| 73) Anthracene-d10            | 17.174 | 188  | 1213512  | 27.651 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.551 | 212  | 1505127  | 30.373 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.527 | 264  | 1687385  | 29.645 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.216  | 88   | 32810    | 10.131 | ng/uL  | 96       |
| 5) Pyridine                   | 3.616  | 79   | 240104   | 28.320 | ng/ul  | 97       |
| 6) Benzaldehyde               | 6.810  | 77   | 34008    | 6.387  | ng/ul  | 94       |
| 8) Phenol                     | 6.869  | 94   | 352676   | 34.405 | ng/ul  | 100      |
| 10) Bis(2-Chloroethyl)ether   | 7.081  | 93   | 248564   | 32.400 | ng/ul  | 97       |
| 12) 2-Chlorophenol            | 7.222  | 128  | 252242   | 34.892 | ng/ul  | 98       |
| 13) 2-Methylphenol            | 8.092  | 108  | 252678   | 33.966 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.157  | 45   | 282623   | 31.699 | ng/ul  | 98       |
| 16) Acetophenone              | 8.457  | 105  | 428380   | 33.001 | ng/ul  | 96       |
| 17) N-Nitroso-di-n-propyla... | 8.445  | 70   | 226081   | 34.376 | ng/ul  | 97       |
| 18) 4-Methylphenol            | 8.416  | 108  | 282056   | 34.177 | ng/ul  | 97       |
| 19) Hexachloroethane          | 8.704  | 117  | 111516   | 33.122 | ng/ul  | 99       |
| 22) Nitrobenzene              | 8.834  | 77   | 340434   | 31.823 | ng/ul  | 98       |
| 23) Isophorone                | 9.357  | 82   | 634616   | 33.086 | ng/ul  | 98       |
| 25) 2-Nitrophenol             | 9.539  | 139  | 158809   | 35.514 | ng/ul  | 95       |
| 26) 2,4-Dimethylphenol        | 9.598  | 107  | 276578   | 28.456 | ng/ul  | 97       |
| 27) Bis(2-Chloroethoxy)met... | 9.822  | 93   | 351726   | 32.296 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.069 | 162  | 311751   | 35.898 | ng/ul  | 99       |
| 30) Naphthalene               | 10.457 | 128  | 810583   | 32.651 | ng/ul  | 98       |
| 32) 4-Chloroaniline           | 10.569 | 127  | 198595   | 19.307 | ng/ul  | 99       |
| 33) Hexachlorobutadiene       | 10.739 | 225  | 248056   | 33.890 | ng/ul  | 98       |
| 34) Caprolactam               | 11.363 | 113  | 96645m   | 34.519 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.710 | 107  | 342837   | 36.183 | ng/ul  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102524\  
 Data File : BP022610.D  
 Acq On : 25 Oct 2024 13:55  
 Operator : RC/JU  
 Sample : P4435-03MSD  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHK68MSD

Manual IntegrationsAPPROVED

Quant Time: Oct 25 14:52:28 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 22 05:59:59 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/26/2024  
 Supervised By :mohammad ahmed 10/29/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.069 | 142  | 628693   | 34.499 | ng/ul  | 98       |
| 37) 1-Methylnaphthalene       | 12.292 | 142  | 613296   | 32.972 | ng/ul  | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.439 | 216  | 473220   | 33.590 | ng/ul  | 96       |
| 40) Hexachlorocyclopentadiene | 12.416 | 237  | 177236   | 20.649 | ng/ul  | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.692 | 196  | 311788   | 35.184 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol     | 12.775 | 196  | 336827   | 35.697 | ng/ul  | 98       |
| 43) 1,1'-Biphenyl             | 13.092 | 154  | 905753   | 32.981 | ng/ul  | 100      |
| 44) 2-Chloronaphthalene       | 13.133 | 162  | 738556   | 32.852 | ng/ul  | 97       |
| 45) 2-Nitroaniline            | 13.351 | 65   | 222183   | 33.098 | ng/ul  | 90       |
| 47) Dimethylphthalate         | 13.733 | 163  | 987114   | 32.724 | ng/ul  | 100      |
| 48) 2,6-Dinitrotoluene        | 13.851 | 165  | 212487   | 37.242 | ng/ul  | 94       |
| 50) Acenaphthylene            | 13.986 | 152  | 1086347  | 32.556 | ng/ul  | 100      |
| 51) 3-Nitroaniline            | 14.186 | 138  | 174541   | 32.571 | ng/ul  | 93       |
| 52) Acenaphthene              | 14.333 | 153  | 770305   | 32.645 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.410 | 184  | 143099   | 34.089 | ng/ul  | 97       |
| 55) 4-Nitrophenol             | 14.521 | 109  | 189593   | 33.637 | ng/ul  | 93       |
| 56) Dibenzofuran              | 14.669 | 168  | 1188333  | 33.527 | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene        | 14.657 | 165  | 323077   | 36.440 | ng/ul  | 95       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.904 | 232  | 318225   | 36.132 | ng/ul  | 99       |
| 59) Diethylphthalate          | 15.116 | 149  | 958764   | 33.130 | ng/ul  | 99       |
| 61) Fluorene                  | 15.333 | 166  | 950556   | 32.914 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.321 | 204  | 544259   | 32.955 | ng/ul  | 98       |
| 63) 4-Nitroaniline            | 15.363 | 138  | 202653   | 37.582 | ng/ul  | 94       |
| 66) 4,6-Dinitro-2-methylph... | 15.445 | 198  | 221768   | 33.555 | ng/ul# | 95       |
| 67) N-Nitrosodiphenylamine    | 15.545 | 169  | 840980   | 33.662 | ng/ul  | 100      |
| 68) 4-Bromophenyl-phenylether | 16.239 | 248  | 389668   | 34.636 | ng/ul  | 96       |
| 69) Hexachlorobenzene         | 16.351 | 284  | 483160   | 34.890 | ng/ul  | 98       |
| 70) Atrazine                  | 16.527 | 200  | 368195   | 35.455 | ng/ul  | 95       |
| 71) Pentachlorophenol         | 16.715 | 266  | 303064   | 34.036 | ng/ul  | 99       |
| 72) Phenanthrene              | 17.110 | 178  | 1665171  | 33.362 | ng/ul  | 99       |
| 74) Anthracene                | 17.210 | 178  | 1613114  | 32.385 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.057 | 216  | 459338   | 32.336 | ng/uL  | 97       |
| 76) Pentachlorobenzene        | 14.592 | 250  | 510090   | 32.483 | ng/uL  | 99       |
| 77) Carbazole                 | 17.486 | 167  | 1430332  | 32.453 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.068 | 149  | 1605417  | 32.920 | ng/ul  | 100      |
| 80) Fluoranthene              | 19.198 | 202  | 2109331  | 37.026 | ng/ul  | 98       |
| 82) Pyrene                    | 19.580 | 202  | 2228590  | 36.498 | ng/ul  | 98       |
| 83) Butylbenzylphthalate      | 20.533 | 149  | 763220   | 35.902 | ng/ul  | 96       |
| 84) 3,3'-Dichlorobenzidine    | 21.403 | 252  | 697680   | 31.233 | ng/ul  | 98       |
| 85) Benzo(a)anthracene        | 21.480 | 228  | 2218963  | 33.778 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.409 | 149  | 1086247  | 35.598 | ng/ul  | 99       |
| 87) Chrysene                  | 21.551 | 228  | 2023180  | 33.214 | ng/ul  | 100      |
| 89) Di-n-octyl phthalate      | 22.662 | 149  | 1787327  | 35.255 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.739 | 252  | 2396427  | 35.720 | ng/ul# | 98       |
| 91) Benzo(k)fluoranthene      | 23.797 | 252  | 2285441  | 34.794 | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 24.603 | 252  | 2099244  | 33.998 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.480 | 276  | 2831559  | 34.365 | ng/ul  | 97       |
| 95) Dibenzo(a,h)anthracene    | 28.538 | 278  | 2280347  | 34.177 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 29.609 | 276  | 2172386m | 33.896 | ng/ul  |          |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

BHK68MSD

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 10/26/2024

Supervised By :mohammad ahmed 10/29/2024

