

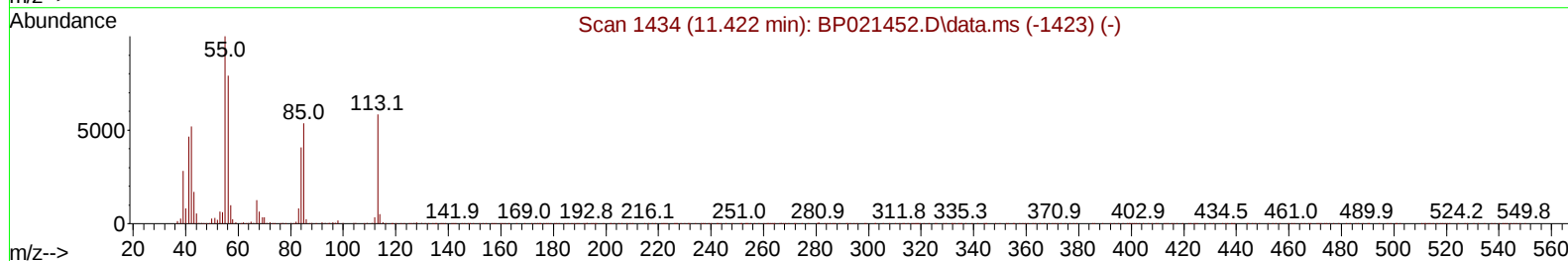
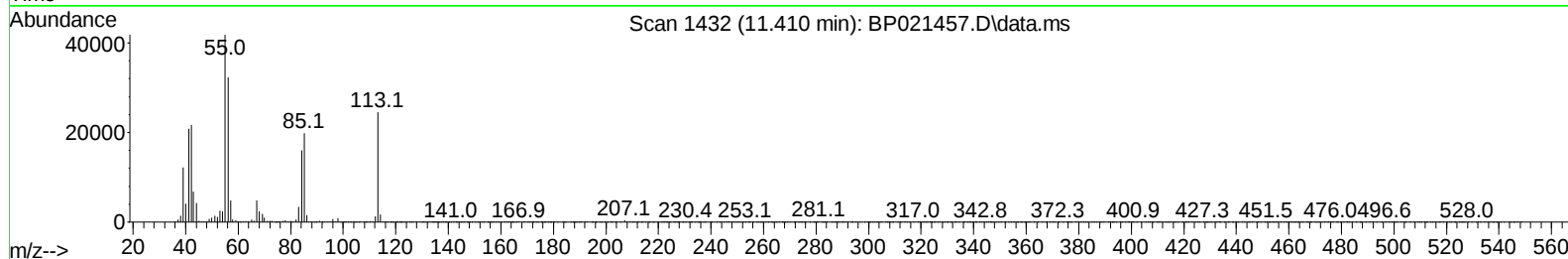
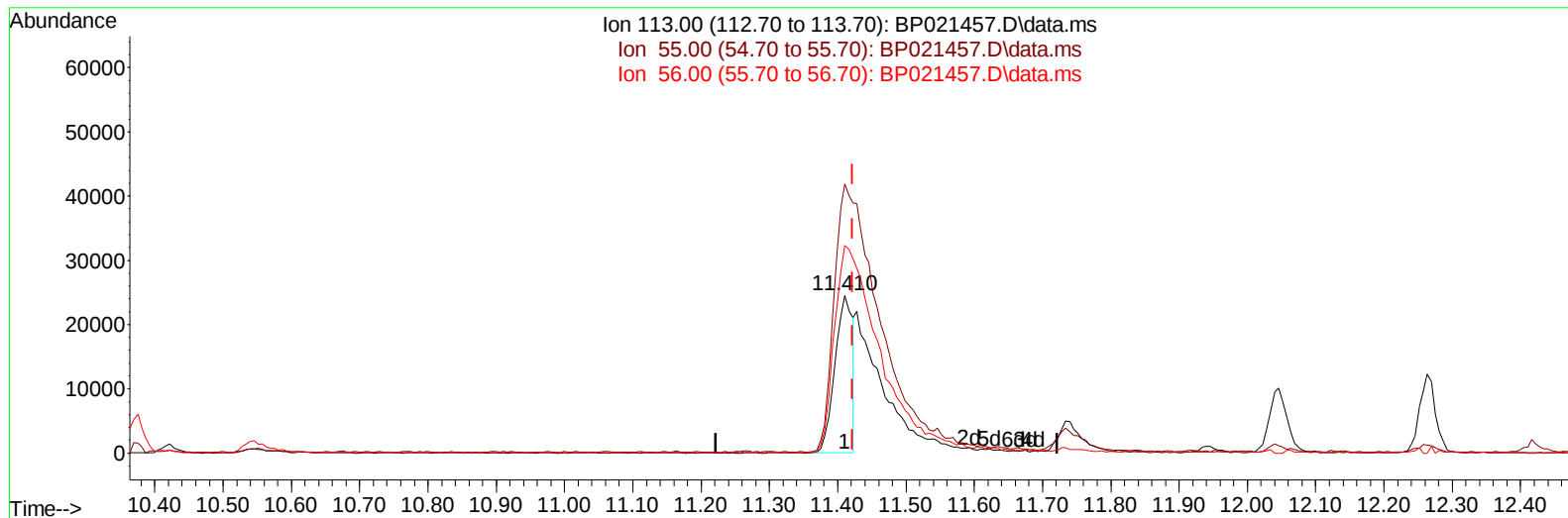
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080824\  
Data File : BP021457.D  
Acq On : 08 Aug 2024 20:57  
Operator : RC/JU  
Sample : PB162539BS  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS539

Manual IntegrationsAPPROVED

Quant Time: Aug 09 02:00:23 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 08 18:23:21 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/09/2024  
Supervised By :mohammad ahmed 08/12/2024



TIC: BP021457.D\data.ms

(34) Caprolactam

11.410min (-0.012) 11.71 ng/ul

response 44232

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 175.30 | 170.42 |
| 56.00  | 131.90 | 131.50 |
| 0.00   | 0.00   | 0.00   |

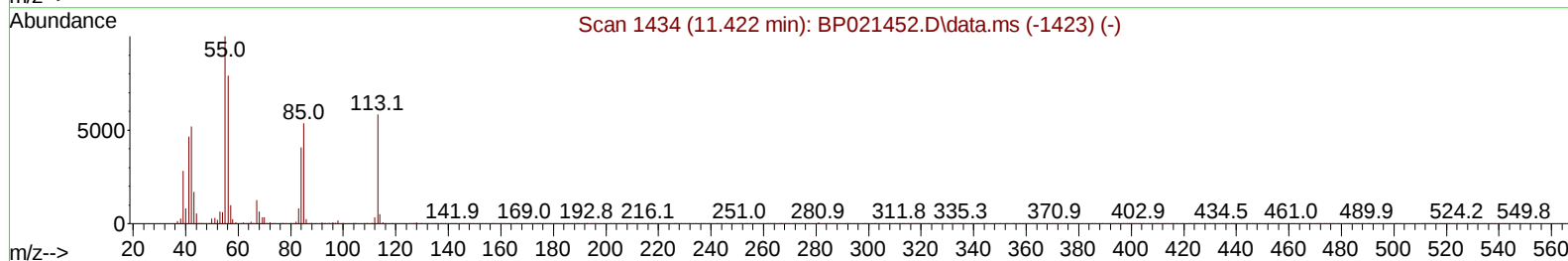
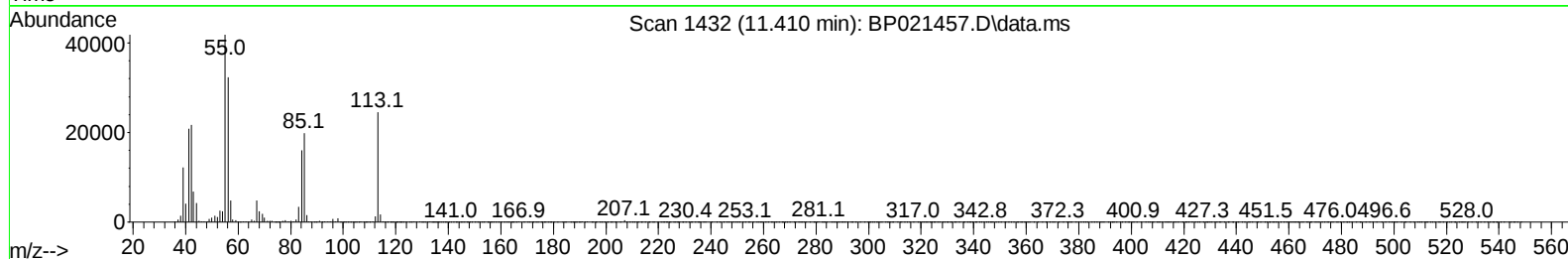
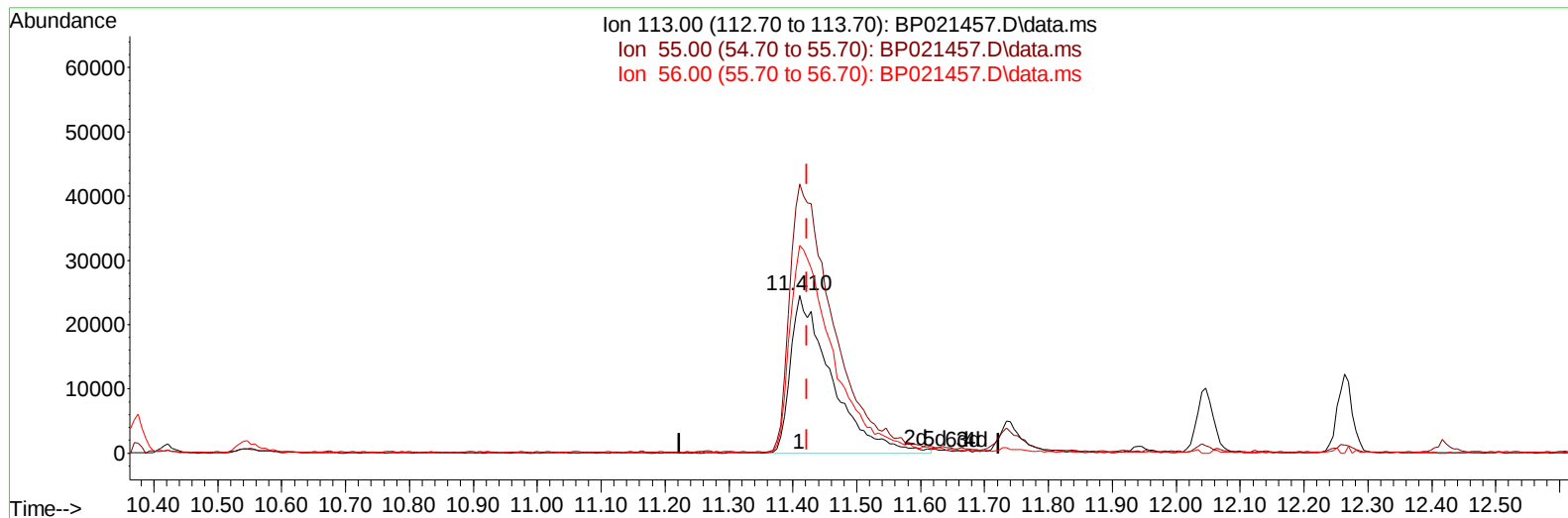
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080824\  
Data File : BP021457.D  
Acq On : 08 Aug 2024 20:57  
Operator : RC/JU  
Sample : PB162539BS  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS539

Manual IntegrationsAPPROVED

Quant Time: Aug 09 02:00:23 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 08 18:23:21 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/09/2024  
Supervised By :mohammad ahmed 08/12/2024



TIC: BP021457.D\data.ms

(34) Caprolactam

11.410min (-0.012) 29.00 ng/ul m

response 109581

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 175.30 | 170.42 |
| 56.00  | 131.90 | 131.50 |
| 0.00   | 0.00   | 0.00   |

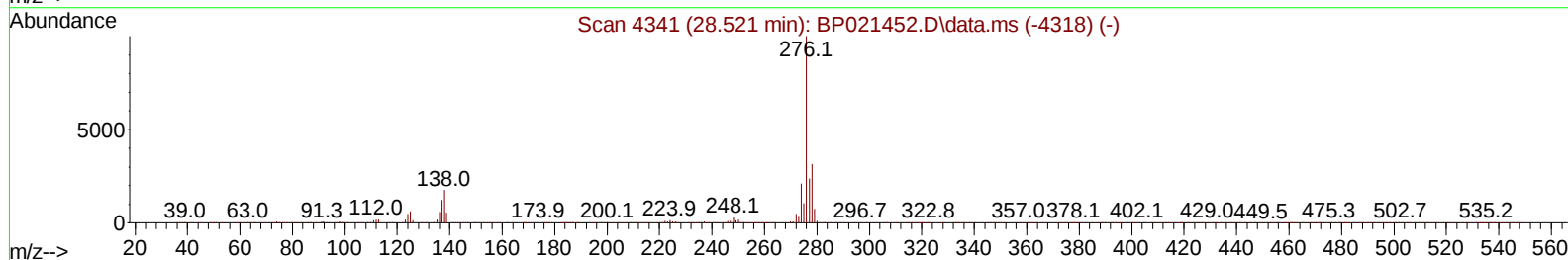
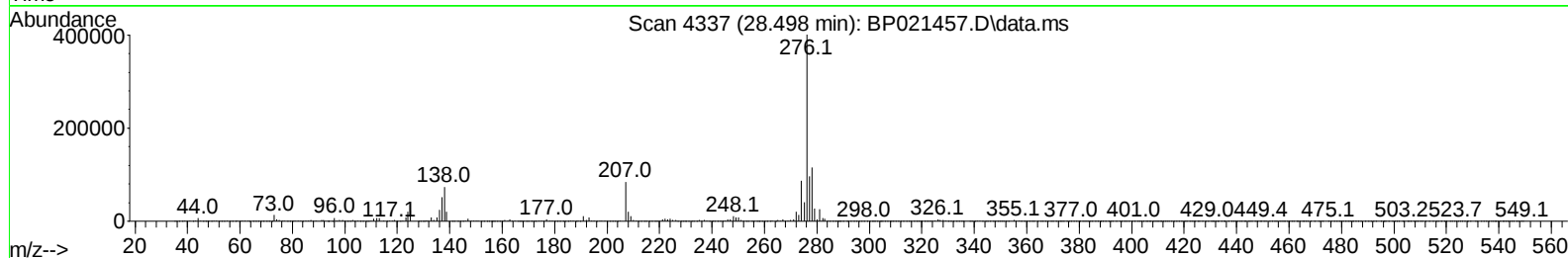
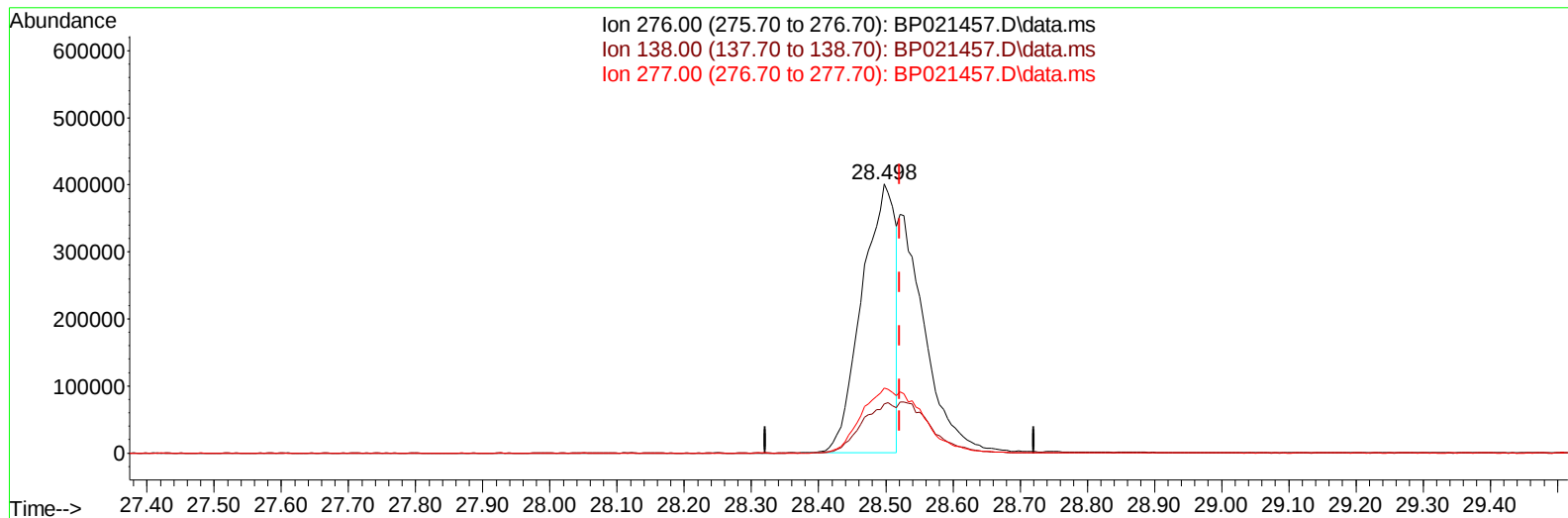
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080824\  
Data File : BP021457.D  
Acq On : 08 Aug 2024 20:57  
Operator : RC/JU  
Sample : PB162539BS  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS539

Manual IntegrationsAPPROVED

Quant Time: Aug 09 02:00:23 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 08 18:23:21 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/09/2024  
Supervised By :mohammad ahmed 08/12/2024



TIC: BP021457.D\data.ms

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

**28.498min (-0.023) 15.81 ng/u1**

**response 1381524**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 17.90  | 23.29# |
| 277.00 | 24.00  | 24.18  |
| 0.00   | 0.00   | 0.00   |

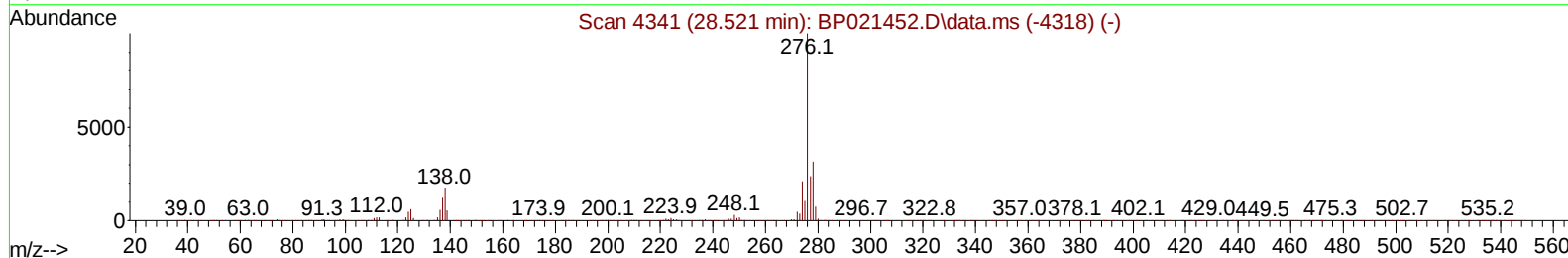
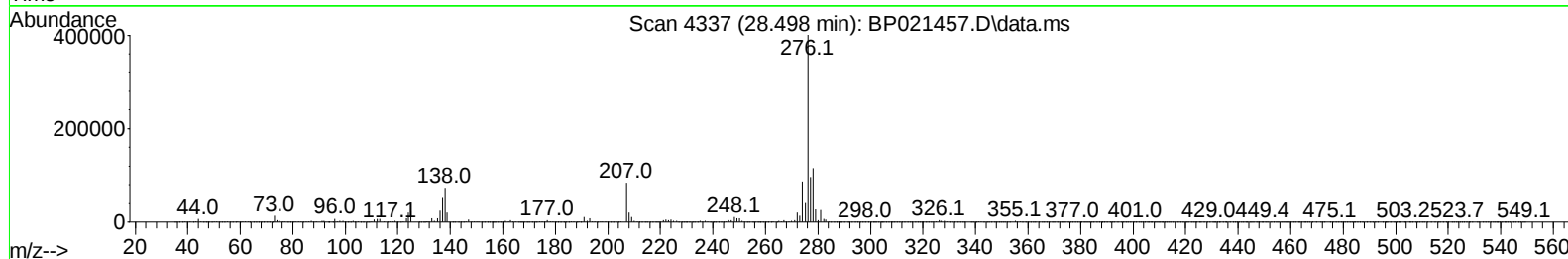
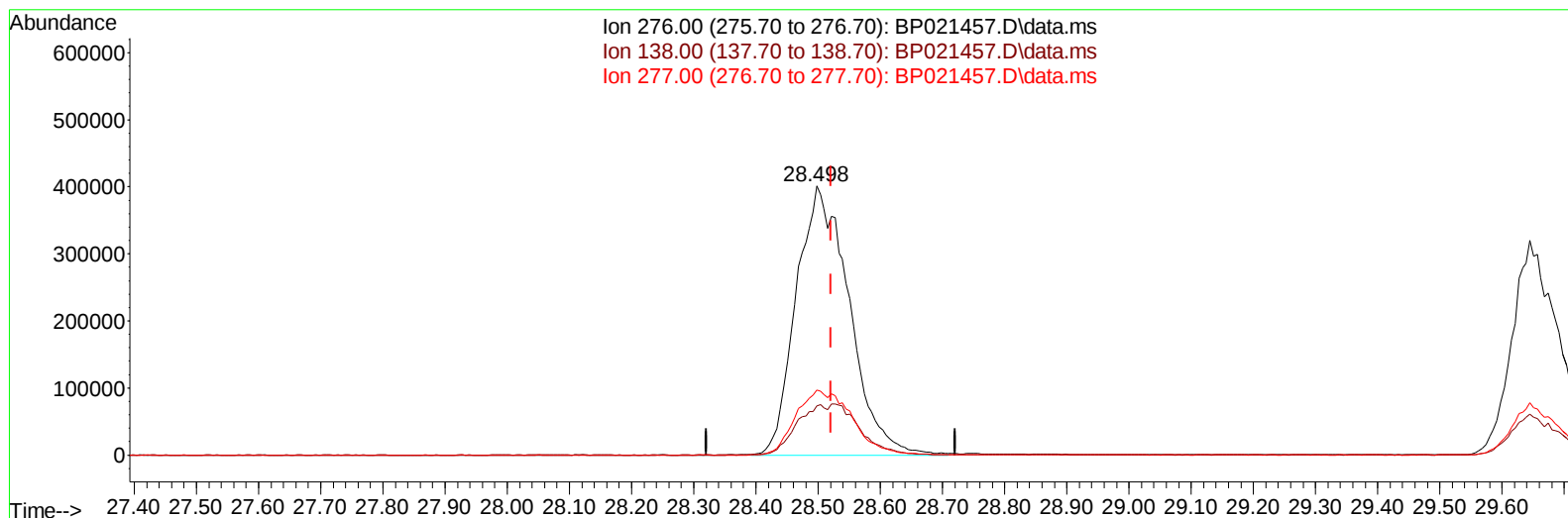
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080824\  
Data File : BP021457.D  
Acq On : 08 Aug 2024 20:57  
Operator : RC/JU  
Sample : PB162539BS  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS539

Manual IntegrationsAPPROVED

Quant Time: Aug 09 02:00:23 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 08 18:23:21 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/09/2024  
Supervised By :mohammad ahmed 08/12/2024



TIC: BP021457.D\data.ms

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

28.498min (-0.023) 27.11 ng/ul m

response 2369185

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 17.90  | 18.32  |
| 277.00 | 24.00  | 24.18  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080824\  
 Data File : BP021457.D  
 Acq On : 08 Aug 2024 20:57  
 Operator : RC/JU  
 Sample : PB162539BS  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS539

Manual IntegrationsAPPROVED

Quant Time: Aug 09 02:51:38 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 08 18:23:21 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/09/2024  
 Supervised By :mohammad ahmed 08/12/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.599  | 152  | 187961   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.375 | 136  | 774284   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.251 | 164  | 429083   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.075 | 188  | 938704   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.510 | 240  | 1051902  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.769 | 264  | 1182718  | 20.000 | ng/ul  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.152  | 96   | 25555    | 6.187  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.570  | 84   | 261314   | 21.853 | ng/ul  | -0.01    |
| 7) Phenol-d5                  | 6.822  | 99   | 393637   | 25.688 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.952  | 67   | 241793   | 26.037 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.152  | 132  | 318740   | 25.245 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.340  | 113  | 272561   | 21.416 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.769  | 128  | 153304   | 25.491 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.481  | 143  | 179624   | 26.096 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.046 | 165  | 323546   | 25.995 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.546 | 131  | 400551   | 24.292 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.663 | 166  | 862798   | 27.661 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.946 | 160  | 952805   | 27.204 | ng/ul  | 0.01     |
| 54) 4-Nitrophenol-d4          | 14.581 | 143  | 154476   | 34.808 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.269 | 176  | 734004   | 28.684 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.457 | 200  | 150412   | 26.483 | ng/ul  | 0.01     |
| 73) Anthracene-d10            | 17.175 | 188  | 1132201  | 27.156 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.551 | 212  | 1416872  | 21.787 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.551 | 264  | 1644098  | 28.185 | ng/ul  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.181  | 88   | 43982    | 9.682  | ng/uL  | 91       |
| 5) Pyridine                   | 3.587  | 79   | 257414   | 20.870 | ng/ul  | 96       |
| 6) Benzaldehyde               | 6.787  | 77   | 208478   | 26.986 | ng/ul  | 96       |
| 8) Phenol                     | 6.852  | 94   | 415915   | 26.822 | ng/ul  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 7.046  | 93   | 326105   | 25.458 | ng/ul  | 98       |
| 12) 2-Chlorophenol            | 7.187  | 128  | 335022   | 26.475 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.069  | 108  | 313489   | 26.141 | ng/ul  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.122  | 45   | 494503   | 26.720 | ng/ul# | 95       |
| 16) Acetophenone              | 8.434  | 105  | 462708   | 23.556 | ng/ul  | 91       |
| 17) N-Nitroso-di-n-propyla... | 8.405  | 70   | 233376   | 22.675 | ng/ul  | 99       |
| 18) 4-Methylphenol            | 8.405  | 108  | 309478   | 24.089 | ng/ul  | 94       |
| 19) Hexachloroethane          | 8.640  | 117  | 122953   | 21.622 | ng/ul  | 99       |
| 22) Nitrobenzene              | 8.811  | 77   | 398278   | 27.646 | ng/ul  | 97       |
| 23) Isophorone                | 9.322  | 82   | 735142   | 25.668 | ng/ul  | 98       |
| 25) 2-Nitrophenol             | 9.516  | 139  | 193398   | 26.940 | ng/ul  | 100      |
| 26) 2,4-Dimethylphenol        | 9.587  | 107  | 309735   | 21.880 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 9.799  | 93   | 443128   | 25.454 | ng/ul  | 98       |
| 29) 2,4-Dichlorophenol        | 10.069 | 162  | 315958   | 25.758 | ng/ul  | 99       |
| 30) Naphthalene               | 10.416 | 128  | 1074865  | 26.391 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.569 | 127  | 378462   | 24.036 | ng/ul  | 100      |
| 33) Hexachlorobutadiene       | 10.675 | 225  | 228731   | 24.869 | ng/ul  | 98       |
| 34) Caprolactam               | 11.410 | 113  | 109581m  | 28.999 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.734 | 107  | 374667   | 28.100 | ng/ul  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080824\  
 Data File : BP021457.D  
 Acq On : 08 Aug 2024 20:57  
 Operator : RC/JU  
 Sample : PB162539BS  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS539

Manual IntegrationsAPPROVED

Quant Time: Aug 09 02:51:38 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 08 18:23:21 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/09/2024  
 Supervised By :mohammad ahmed 08/12/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.046 | 142  | 736526   | 26.345 | ng/ul | 99       |
| 37) 1-Methylnaphthalene       | 12.263 | 142  | 736640   | 25.619 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.422 | 216  | 428754   | 30.768 | ng/ul | 98       |
| 40) Hexachlorocyclopentadiene | 12.363 | 237  | 76562    | 10.503 | ng/ul | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.687 | 196  | 237243   | 26.944 | ng/ul | 98       |
| 42) 2,4,5-Trichlorophenol     | 12.804 | 196  | 272246   | 29.223 | ng/ul | 97       |
| 43) 1,1'-Biphenyl             | 13.063 | 154  | 886778   | 28.279 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.110 | 162  | 699234   | 27.932 | ng/ul | 98       |
| 45) 2-Nitroaniline            | 13.363 | 65   | 194618   | 28.112 | ng/ul | 99       |
| 47) Dimethylphthalate         | 13.716 | 163  | 863599   | 27.286 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 13.863 | 165  | 180411   | 28.512 | ng/ul | 96       |
| 50) Acenaphthylene            | 13.975 | 152  | 1066262  | 26.888 | ng/ul | 99       |
| 51) 3-Nitroaniline            | 14.222 | 138  | 171377   | 31.124 | ng/ul | 93       |
| 52) Acenaphthene              | 14.316 | 153  | 707612   | 26.885 | ng/ul | 99       |
| 53) 2,4-Dinitrophenol         | 14.469 | 184  | 87527    | 26.340 | ng/ul | 96       |
| 55) 4-Nitrophenol             | 14.599 | 109  | 112058   | 31.041 | ng/ul | 97       |
| 56) Dibenzofuran              | 14.663 | 168  | 1012700  | 28.097 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 14.704 | 165  | 267932   | 30.846 | ng/ul | 87       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.922 | 232  | 186865   | 23.424 | ng/ul | 95       |
| 59) Diethylphthalate          | 15.104 | 149  | 868060   | 27.444 | ng/ul | 98       |
| 61) Fluorene                  | 15.322 | 166  | 829998   | 28.220 | ng/ul | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.316 | 204  | 445724   | 27.928 | ng/ul | 97       |
| 63) 4-Nitroaniline            | 15.416 | 138  | 162379   | 35.610 | ng/ul | 100      |
| 66) 4,6-Dinitro-2-methylph... | 15.475 | 198  | 131206   | 21.784 | ng/ul | 99       |
| 67) N-Nitrosodiphenylamine    | 15.551 | 169  | 723745   | 26.477 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.240 | 248  | 291405   | 26.485 | ng/ul | 98       |
| 69) Hexachlorobenzene         | 16.357 | 284  | 353124   | 27.991 | ng/ul | 95       |
| 70) Atrazine                  | 16.545 | 200  | 28034    | 2.814  | ng/ul | 100      |
| 71) Pentachlorophenol         | 16.745 | 266  | 133536   | 19.489 | ng/ul | 98       |
| 72) Phenanthrene              | 17.116 | 178  | 1355859  | 27.843 | ng/ul | 100      |
| 74) Anthracene                | 17.204 | 178  | 1324951  | 26.689 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.028 | 216  | 397053   | 26.427 | ng/uL | 99       |
| 76) Pentachlorobenzene        | 14.581 | 250  | 390000   | 26.123 | ng/uL | 98       |
| 77) Carbazole                 | 17.522 | 167  | 1199294  | 29.552 | ng/ul | 100      |
| 78) Di-n-butylphthalate       | 18.063 | 149  | 1502445  | 26.834 | ng/ul | 99       |
| 80) Fluoranthene              | 19.204 | 202  | 1667772  | 21.236 | ng/ul | 98       |
| 82) Pyrene                    | 19.586 | 202  | 1756726  | 21.981 | ng/ul | 97       |
| 83) Butylbenzylphthalate      | 20.516 | 149  | 706380   | 21.214 | ng/ul | 96       |
| 84) 3,3'-Dichlorobenzidine    | 21.416 | 252  | 627725   | 26.115 | ng/ul | 99       |
| 85) Benzo(a)anthracene        | 21.486 | 228  | 1984143  | 26.927 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.386 | 149  | 1160123  | 24.210 | ng/ul | 99       |
| 87) Chrysene                  | 21.551 | 228  | 1872072  | 27.341 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.616 | 149  | 1975525  | 25.285 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.733 | 252  | 2021111  | 28.160 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.804 | 252  | 1971391  | 27.523 | ng/ul | 99       |
| 93) Benzo(a)pyrene            | 24.622 | 252  | 1798084  | 26.511 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 28.498 | 276  | 2369185m | 27.108 | ng/ul |          |
| 95) Dibenzo(a,h)anthracene    | 28.551 | 278  | 1948955  | 26.929 | ng/ul | 99       |
| 96) Benzo(g,h,i)perylene      | 29.645 | 276  | 1825788  | 25.651 | ng/ul | 98       |

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

**Instrument :**  
BNA\_P  
**ClientSampleId :**  
SLCS539

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 08/09/2024  
Supervised By :mohammad ahmed 08/12/2024

