

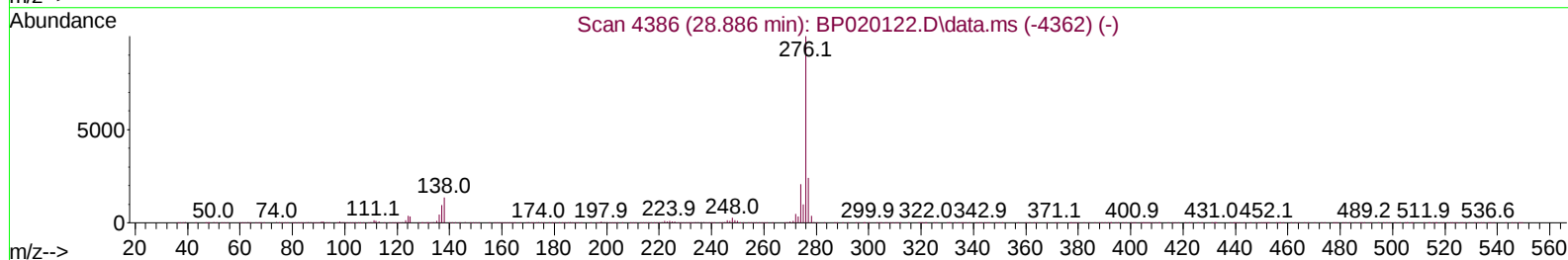
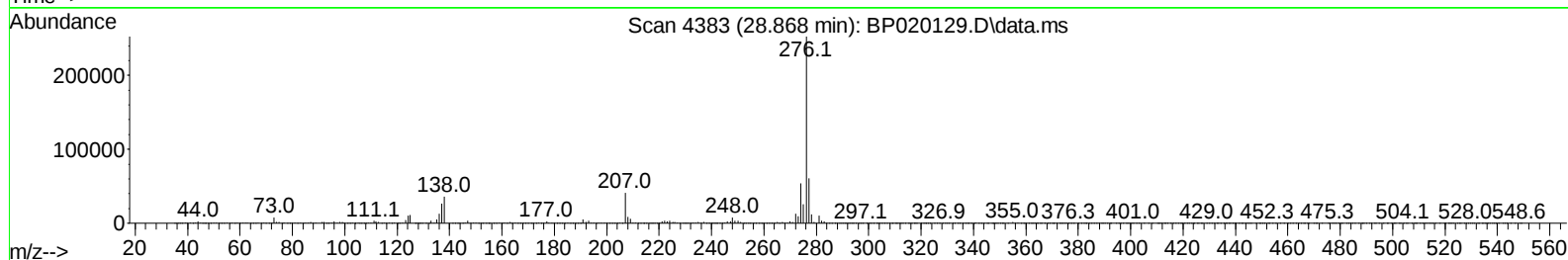
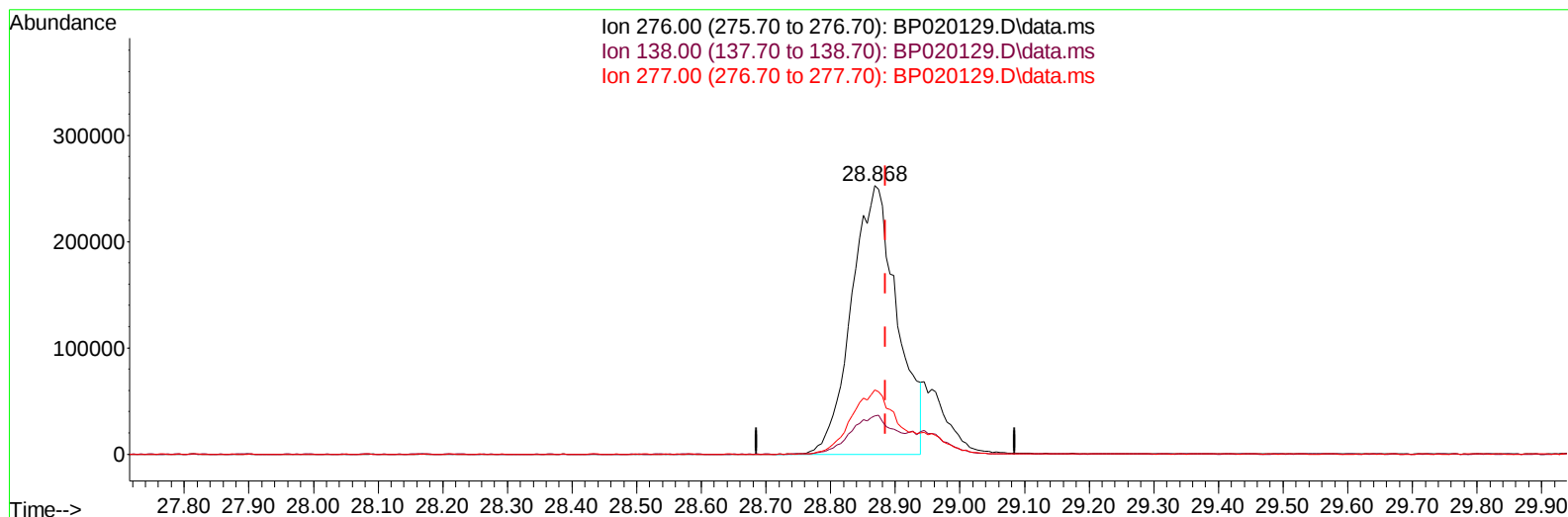
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP043024\  
 Data File : BP020129.D  
 Acq On : 30 Apr 2024 21:48  
 Operator : MA/JU  
 Sample : P2220-05MSD 10X  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 E0813MSD

Manual IntegrationsAPPROVED

Quant Time: May 01 02:03:25 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 01 00:58:38 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/03/2024  
 Supervised By :mohammad ahmed 05/03/2024



TIC: BP020129.D\data.ms

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

**28.868min (-0.017) 46.79 ng/u1**

**response 1234604**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 12.40  | 14.25  |
| 277.00 | 24.50  | 23.93  |
| 0.00   | 0.00   | 0.00   |

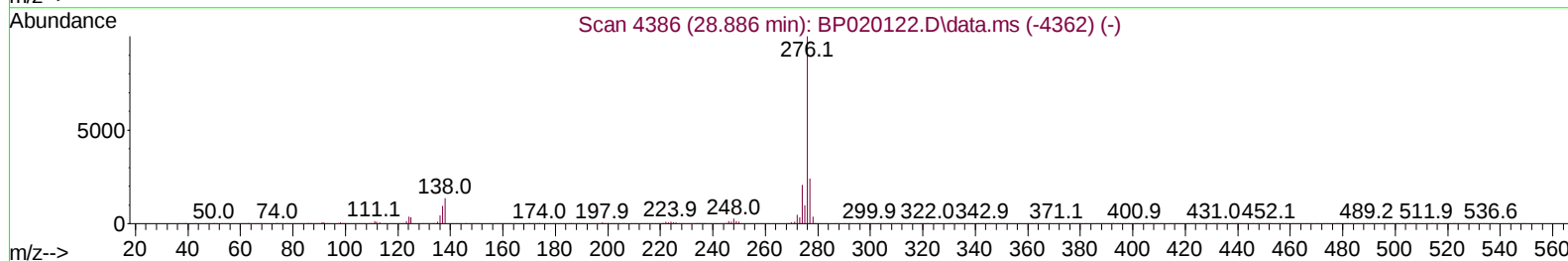
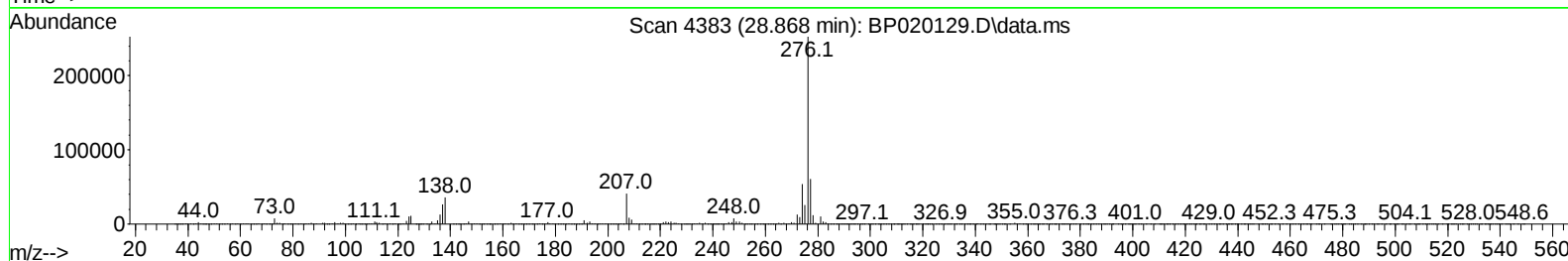
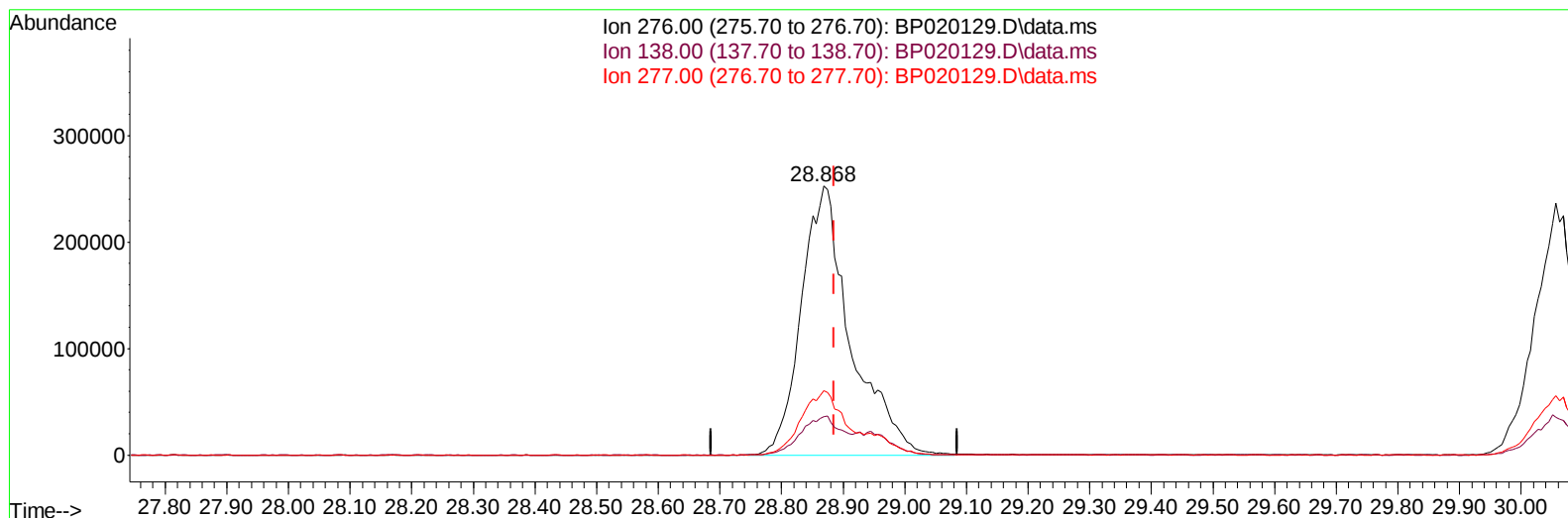
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP043024\  
Data File : BP020129.D  
Acq On : 30 Apr 2024 21:48  
Operator : MA/JU  
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Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
E0813MSD

Manual IntegrationsAPPROVED

Quant Time: May 01 02:03:25 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042424.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed May 01 00:58:38 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/03/2024  
Supervised By :mohammad ahmed 05/03/2024



TIC: BP020129.D\data.ms

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

**28.868min (-0.017) 53.29 ng/ul m**

**response 1406148**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 12.40  | 14.25  |
| 277.00 | 24.50  | 23.93  |
| 0.00   | 0.00   | 0.00   |

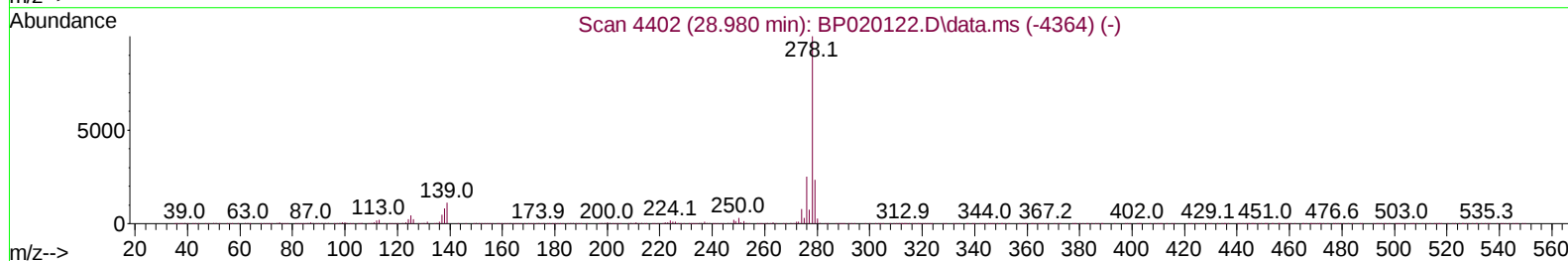
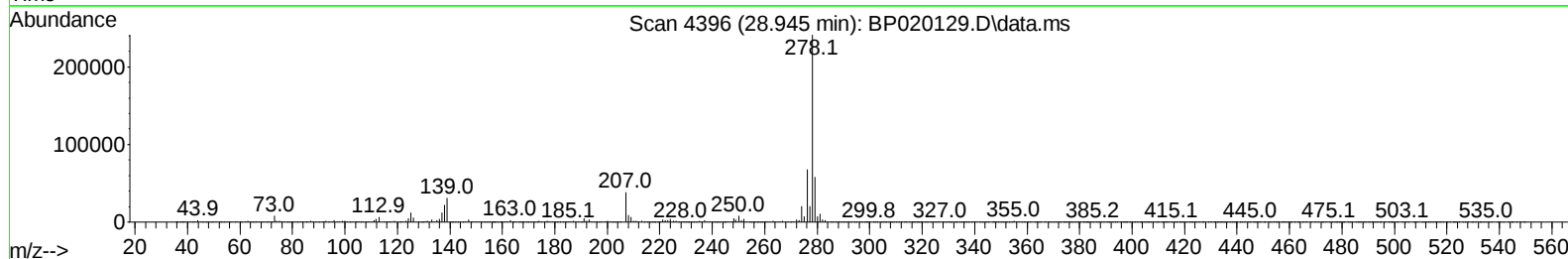
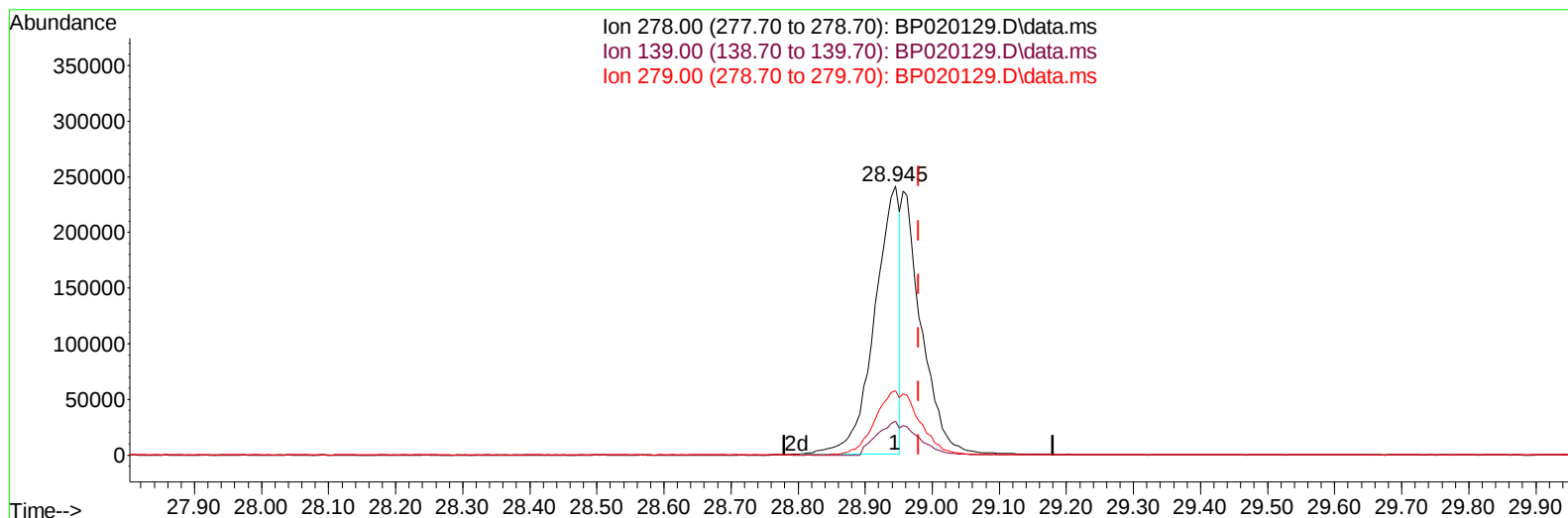
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP043024\  
 Data File : BP020129.D  
 Acq On : 30 Apr 2024 21:48  
 Operator : MA/JU  
 Sample : P2220-05MSD 10X  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 E0813MSD

Manual IntegrationsAPPROVED

Quant Time: May 01 02:03:25 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 01 00:58:38 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/03/2024  
 Supervised By :mohammad ahmed 05/03/2024



TIC: BP020129.D\data.ms

**(95) Dibenzo(a,h)anthracene**

**28.945min (-0.035) 28.56 ng/u1**

**response 622071**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 278.00 | 100.00 | 100.00 |
| 139.00 | 11.80  | 12.74  |
| 279.00 | 24.30  | 23.93  |
| 0.00   | 0.00   | 0.00   |

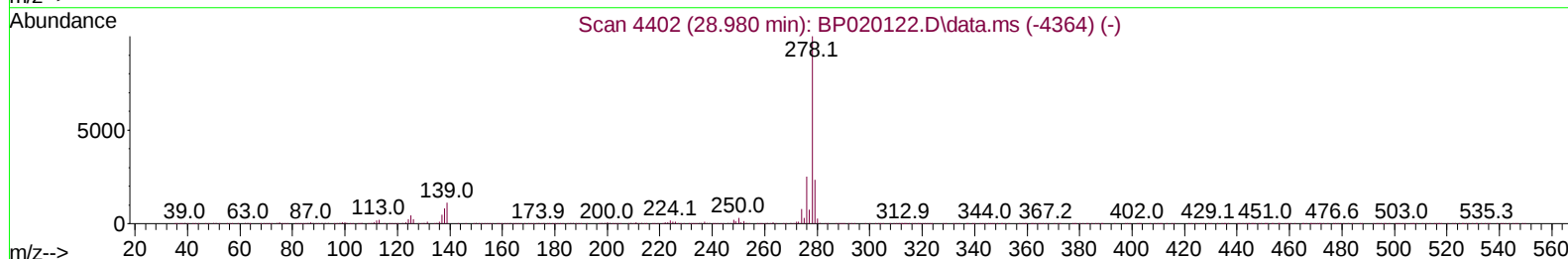
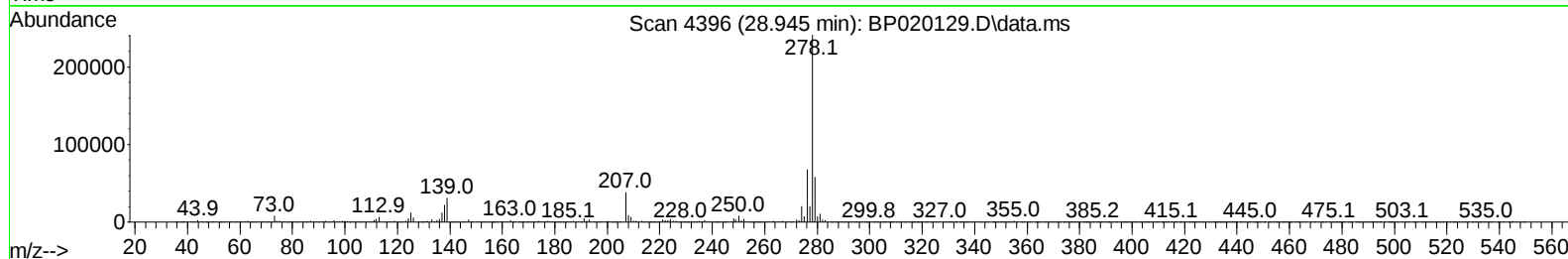
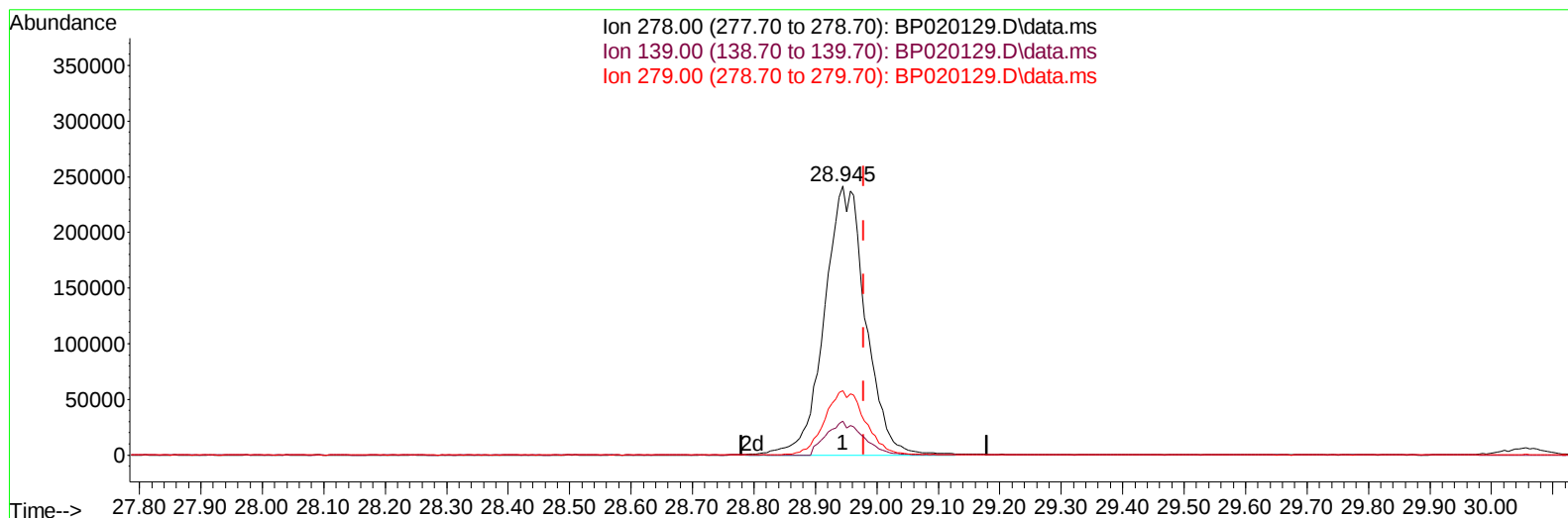
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP043024\  
 Data File : BP020129.D  
 Acq On : 30 Apr 2024 21:48  
 Operator : MA/JU  
 Sample : P2220-05MSD 10X  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 E0813MSD

Manual IntegrationsAPPROVED

Quant Time: May 01 02:03:25 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 01 00:58:38 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/03/2024  
 Supervised By :mohammad ahmed 05/03/2024



TIC: BP020129.D\data.ms

**(95) Dibenzo(a,h)anthracene**

**28.945min (-0.035) 51.71 ng/ul m**

**response 1126296**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 278.00 | 100.00 | 100.00 |
| 139.00 | 11.80  | 12.74  |
| 279.00 | 24.30  | 23.93  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP043024\  
 Data File : BP020129.D  
 Acq On : 30 Apr 2024 21:48  
 Operator : MA/JU  
 Sample : P2220-05MSD 10X  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 E0813MSD

Manual IntegrationsAPPROVED

Quant Time: May 01 02:05:14 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 01 00:58:38 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/03/2024  
 Supervised By :mohammad ahmed 05/03/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.628  | 152  | 87410    | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.410 | 136  | 371944   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.310 | 164  | 266076   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.151 | 188  | 557820   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.645 | 240  | 384269   | 20.000 | ng/ul  | -0.02    |
| 88) Perylene-d12              | 25.027 | 264  | 385052   | 20.000 | ng/ul  | -0.02    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.223  | 96   | 5827     | 2.784  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.628  | 84   | 77255    | 12.529 | ng/ul  | 0.01     |
| 7) Phenol-d5                  | 6.822  | 99   | 68485    | 9.240  | ng/ul  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.987  | 67   | 213548   | 44.724 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.169  | 132  | 172509   | 33.149 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.340  | 113  | 132029   | 23.397 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.793  | 128  | 122360   | 42.839 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.505  | 143  | 134388   | 42.395 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.040 | 165  | 237828   | 39.867 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.563 | 131  | 310591   | 39.925 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.728 | 166  | 902516   | 47.491 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.998 | 160  | 958322   | 44.243 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.569 | 143  | 26233    | 8.492  | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.334 | 176  | 746378   | 45.152 | ng/ul  | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.487 | 200  | 163270   | 45.776 | ng/ul  | -0.01    |
| 73) Anthracene-d10            | 17.257 | 188  | 1188102  | 47.972 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.651 | 212  | 1283639  | 62.789 | ng/ul  | -0.01    |
| 92) Benzo(a)pyrene-d12        | 24.798 | 264  | 932168   | 49.616 | ng/ul  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.252  | 88   | 14355    | 6.101  | ng/ul# | 1        |
| 5) Pyridine                   | 3.646  | 79   | 88636    | 14.016 | ng/ul  | 98       |
| 6) Benzaldehyde               | 6.799  | 77   | 240488   | 56.652 | ng/ul  | 99       |
| 8) Phenol                     | 6.846  | 94   | 76201    | 9.895  | ng/ul  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.075  | 93   | 256720   | 42.278 | ng/ul  | 98       |
| 12) 2-Chlorophenol            | 7.205  | 128  | 178284   | 32.645 | ng/ul  | 96       |
| 13) 2-Methylphenol            | 8.075  | 108  | 147913   | 26.576 | ng/ul  | 95       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.152  | 45   | 357793   | 42.764 | ng/ul  | 99       |
| 16) Acetophenone              | 8.463  | 105  | 439337   | 46.883 | ng/ul  | 100      |
| 17) N-Nitroso-di-n-propyla... | 8.440  | 70   | 250379   | 48.824 | ng/ul# | 99       |
| 18) 4-Methylphenol            | 8.405  | 108  | 142121   | 24.125 | ng/ul  | 96       |
| 19) Hexachloroethane          | 8.681  | 117  | 93634    | 35.481 | ng/ul  | 98       |
| 22) Nitrobenzene              | 8.840  | 77   | 357779   | 41.845 | ng/ul  | 96       |
| 23) Isophorone                | 9.358  | 82   | 697978   | 45.868 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.534  | 139  | 137084   | 40.947 | ng/ul  | 99       |
| 26) 2,4-Dimethylphenol        | 9.599  | 107  | 250230   | 35.583 | ng/ul  | 96       |
| 27) Bis(2-Chloroethoxy)met... | 9.840  | 93   | 379216   | 43.712 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.069 | 162  | 229159   | 39.019 | ng/ul  | 98       |
| 30) Naphthalene               | 10.463 | 128  | 771860   | 40.140 | ng/ul  | 98       |
| 32) 4-Chloroaniline           | 10.587 | 127  | 296783   | 38.578 | ng/ul  | 97       |
| 33) Hexachlorobutadiene       | 10.722 | 225  | 180631   | 36.830 | ng/ul  | 95       |
| 34) Caprolactam               | 11.416 | 113  | 13672    | 6.908  | ng/ul  | 98       |
| 35) 4-Chloro-3-methylphenol   | 11.722 | 107  | 265341   | 39.022 | ng/ul  | 99       |

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 Sample : P2220-05MSD 10X  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 E0813MSD

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 05/03/2024  
 Supervised By :mohammad ahmed 05/03/2024

Quant Time: May 01 02:05:14 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 01 00:58:38 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.087 | 142  | 569523   | 44.227 | ng/ul  | 99       |
| 37) 1-Methylnaphthalene       | 12.310 | 142  | 577801   | 44.652 | ng/ul  | 98       |
| 39) 1,2,4,5-Tetrachloroben... | 12.457 | 216  | 373944   | 41.478 | ng/ul  | 98       |
| 40) Hexachlorocyclopentadiene | 12.416 | 237  | 157684   | 33.096 | ng/ul  | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.710 | 196  | 233332   | 42.699 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol     | 12.793 | 196  | 253169   | 43.129 | ng/ul  | 96       |
| 43) 1,1'-Biphenyl             | 13.116 | 154  | 792709   | 42.399 | ng/ul  | 100      |
| 44) 2-Chloronaphthalene       | 13.163 | 162  | 636852   | 42.076 | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.393 | 65   | 254248   | 48.098 | ng/ul  | 100      |
| 47) Dimethylphthalate         | 13.775 | 163  | 872006   | 44.959 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 13.904 | 165  | 184042   | 46.663 | ng/ul  | 97       |
| 50) Acenaphthylene            | 14.028 | 152  | 1028823  | 42.481 | ng/ul  | 100      |
| 51) 3-Nitroaniline            | 14.246 | 138  | 169037   | 45.704 | ng/ul  | 91       |
| 52) Acenaphthene              | 14.381 | 153  | 665209   | 41.905 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.469 | 184  | 108492   | 43.332 | ng/ul  | 94       |
| 55) 4-Nitrophenol             | 14.587 | 109  | 27272    | 8.345  | ng/ul  | 99       |
| 56) Dibenzofuran              | 14.728 | 168  | 1006119  | 44.758 | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene        | 14.722 | 165  | 270765   | 46.849 | ng/ul  | 95       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.963 | 232  | 240979   | 44.593 | ng/ul# | 99       |
| 59) Diethylphthalate          | 15.175 | 149  | 892000   | 45.608 | ng/ul  | 99       |
| 61) Fluorene                  | 15.393 | 166  | 825141   | 44.205 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.393 | 204  | 450341   | 44.155 | ng/ul  | 98       |
| 63) 4-Nitroaniline            | 15.445 | 138  | 153790   | 46.461 | ng/ul  | 98       |
| 66) 4,6-Dinitro-2-methylph... | 15.498 | 198  | 171434   | 45.466 | ng/ul  | 95       |
| 67) N-Nitrosodiphenylamine    | 15.622 | 169  | 738021   | 50.872 | ng/ul  | 98       |
| 68) 4-Bromophenyl-phenylether | 16.316 | 248  | 295476   | 49.120 | ng/ul  | 94       |
| 69) Hexachlorobenzene         | 16.422 | 284  | 324554   | 53.286 | ng/ul  | 100      |
| 70) Atrazine                  | 16.622 | 200  | 292822   | 49.980 | ng/ul  | 99       |
| 71) Pentachlorophenol         | 16.792 | 266  | 196689   | 45.219 | ng/ul  | 98       |
| 72) Phenanthrene              | 17.192 | 178  | 1317656  | 46.198 | ng/ul  | 99       |
| 74) Anthracene                | 17.292 | 178  | 1347416  | 45.792 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.081 | 216  | 370047   | 45.143 | ng/uL  | 99       |
| 76) Pentachlorobenzene        | 14.634 | 250  | 384002   | 47.938 | ng/uL  | 97       |
| 77) Carbazole                 | 17.587 | 167  | 1160328  | 44.440 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.181 | 149  | 1413404  | 44.193 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.304 | 202  | 1501786  | 63.393 | ng/ul  | 98       |
| 82) Pyrene                    | 19.681 | 202  | 1536945  | 61.821 | ng/ul  | 99       |
| 83) Butylbenzylphthalate      | 20.657 | 149  | 504848   | 50.340 | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.563 | 252  | 384321   | 42.760 | ng/ul  | 98       |
| 85) Benzo(a)anthracene        | 21.627 | 228  | 1247787  | 46.779 | ng/ul  | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.575 | 149  | 659464   | 44.954 | ng/ul  | 97       |
| 87) Chrysene                  | 21.698 | 228  | 1181985  | 47.747 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.886 | 149  | 1044081  | 45.344 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.951 | 252  | 1119759  | 49.064 | ng/ul  | 100      |
| 91) Benzo(k)fluoranthene      | 24.027 | 252  | 1095292  | 47.457 | ng/ul  | 98       |
| 93) Benzo(a)pyrene            | 24.868 | 252  | 937081   | 42.813 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.868 | 276  | 1406148m | 53.289 | ng/ul  |          |
| 95) Dibenzo(a,h)anthracene    | 28.945 | 278  | 1126296m | 51.709 | ng/ul  |          |
| 96) Benzo(g,h,i)perylene      | 30.056 | 276  | 1105631  | 52.108 | ng/ul  | 99       |

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

**Instrument :**  
BNA\_P  
**ClientSampleId :**  
E0813MSD

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 05/03/2024  
Supervised By :mohammad ahmed 05/03/2024

