

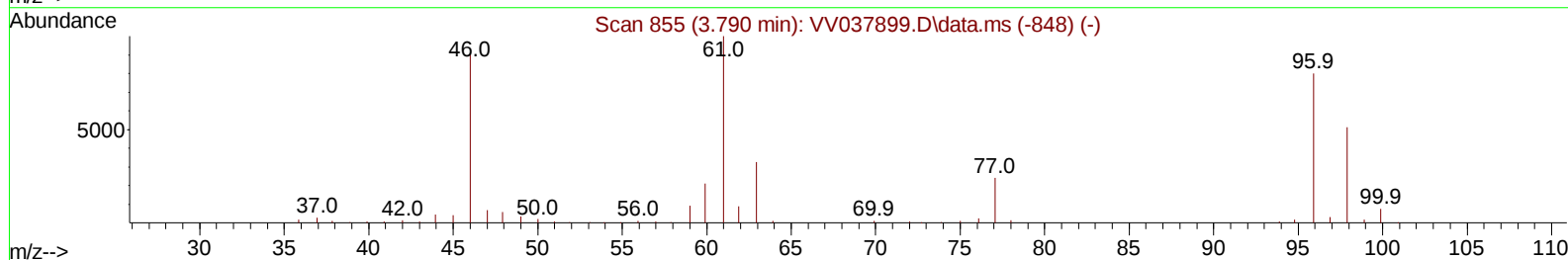
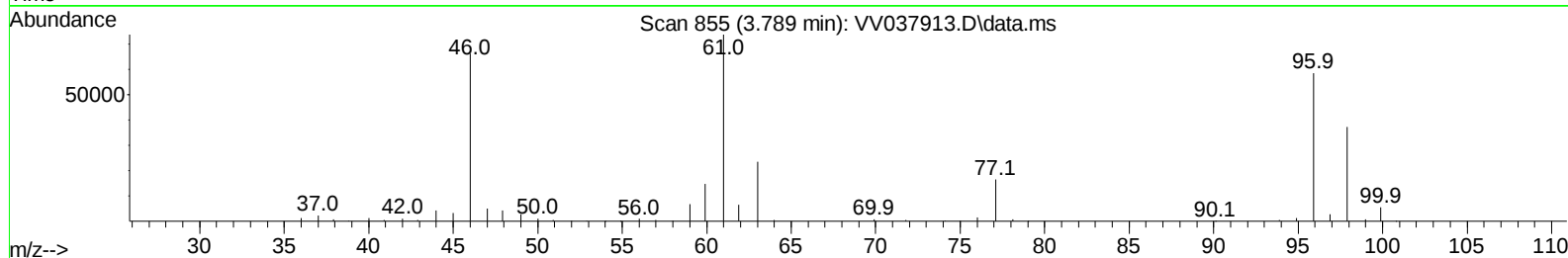
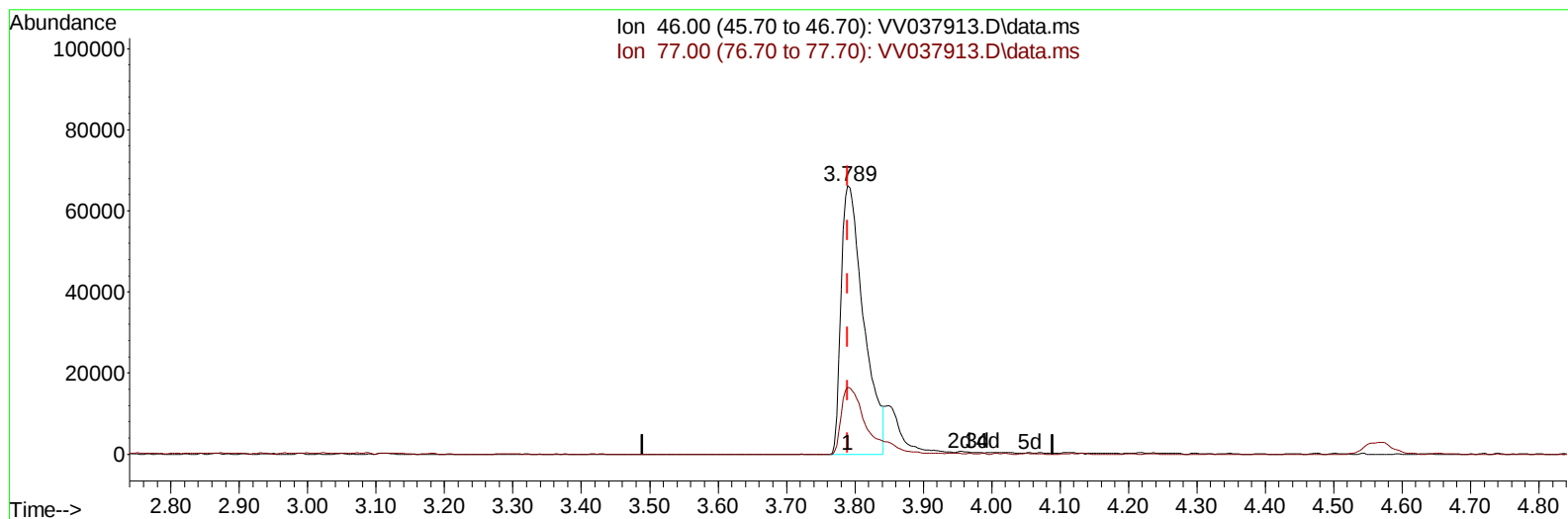
Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\DATA\VV102324\  
Data File : VV037913.D  
Acq On : 23 Oct 2024 10:25  
Operator : SY/MD  
Sample : VSTDCCC050  
Misc : 5mL/MSVOA\_V/WATER  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 01:43:57 2024  
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM092624WMA.M  
Quant Title : VOC Analysis  
QLast Update : Tue Oct 22 05:07:16 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/25/2024  
Supervised By :mohammad ahmed 10/26/2024



TIC: VV037913.D\data.ms

(21) 2-Butanone-d5 (S)

3.789min (-0.000) 102.80 ug/L

response 150932

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 46.00 | 100.00 | 100.00 |
| 77.00 | 25.70  | 28.71  |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

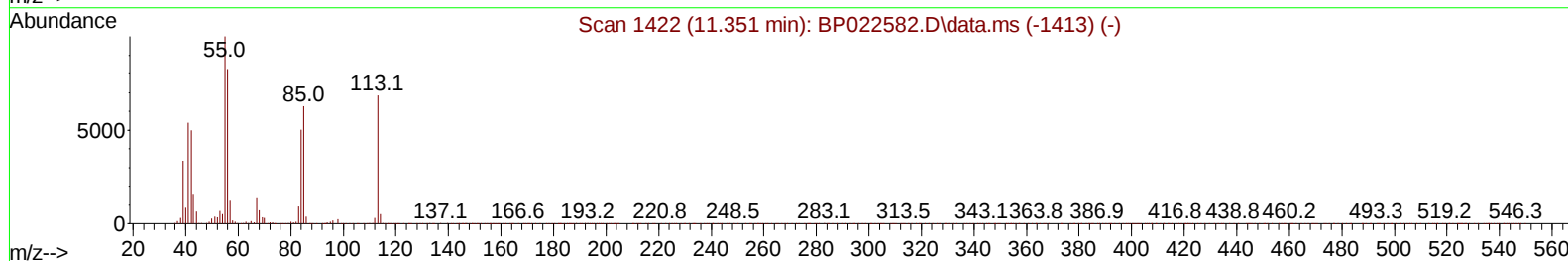
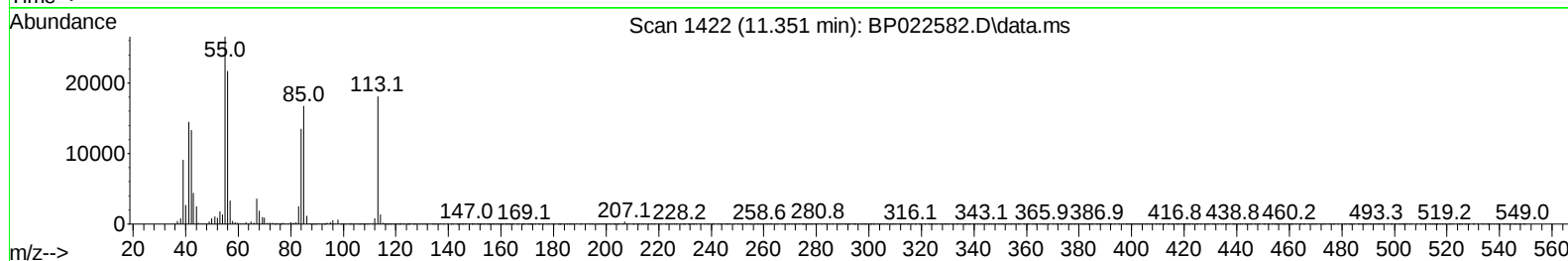
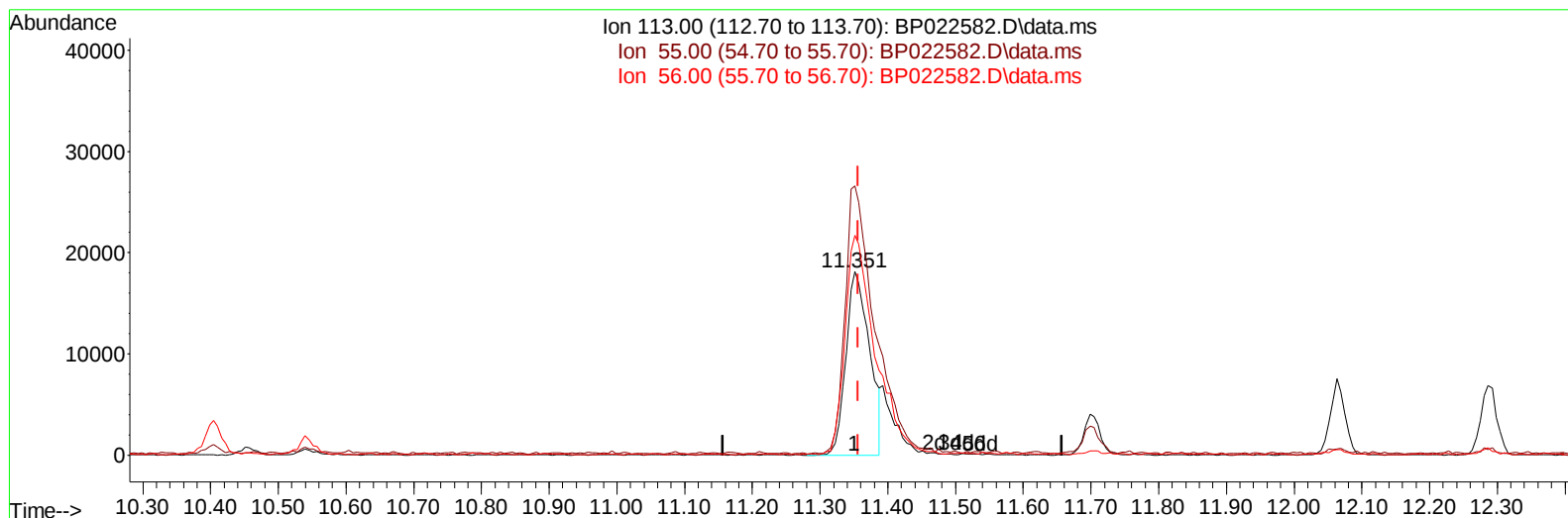
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102424\  
Data File : BP022582.D  
Acq On : 24 Oct 2024 11:00  
Operator : RC/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:23:50 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 10/25/2024  
Supervised By :mohammad ahmed 10/26/2024



TIC: BP022582.D\data.ms

**(34) Caprolactam**

**11.351min (-0.006) 16.18 ng/ul**

**response 43701**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 162.70 | 146.73 |
| 56.00  | 130.00 | 119.74 |
| 0.00   | 0.00   | 0.00   |

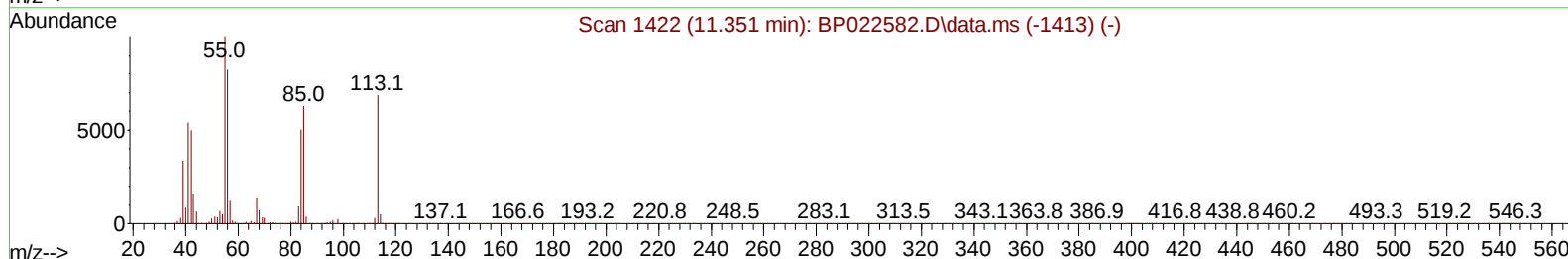
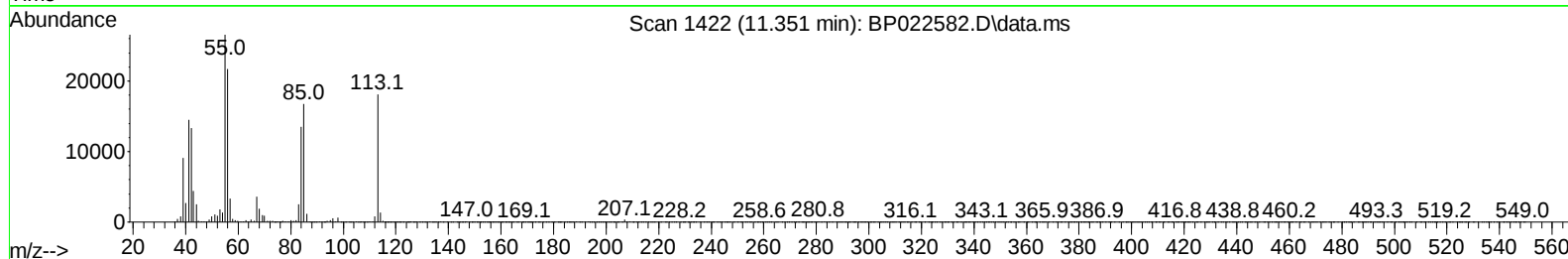
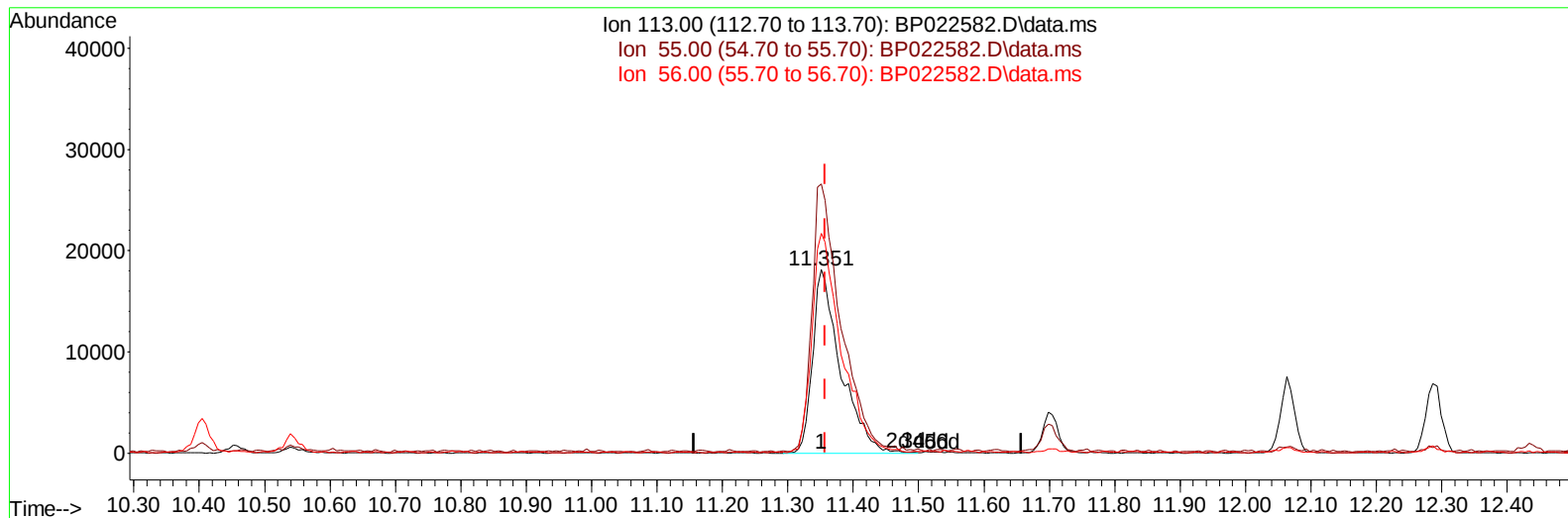
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Data File : BP022582.D  
Acq On : 24 Oct 2024 11:00  
Operator : RC/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:23:50 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 10/25/2024  
Supervised By :mohammad ahmed 10/26/2024



TIC: BP022582.D\data.ms

(34) Caprolactam

11.351min (-0.006) 19.88 ng/ul m

response 53713

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 162.70 | 146.73 |
| 56.00  | 130.00 | 119.74 |
| 0.00   | 0.00   | 0.00   |

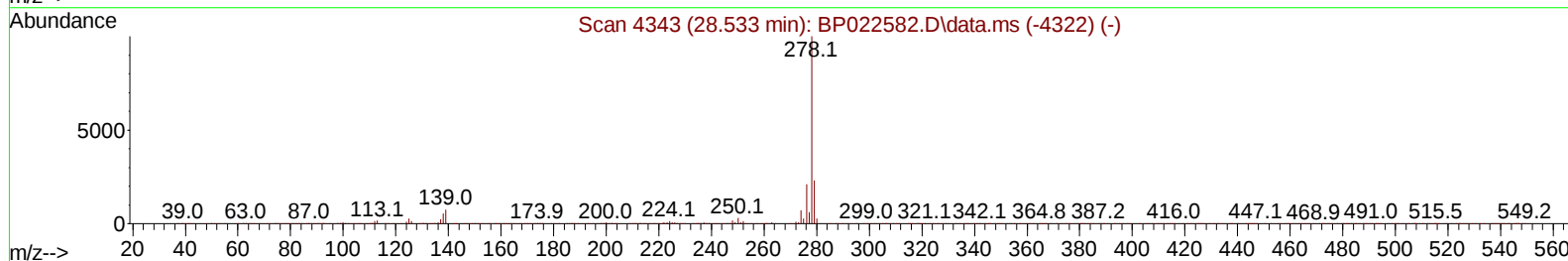
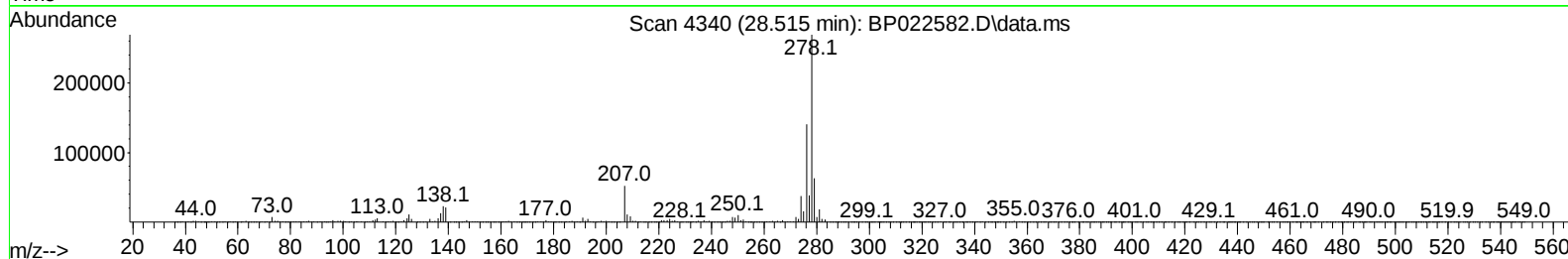
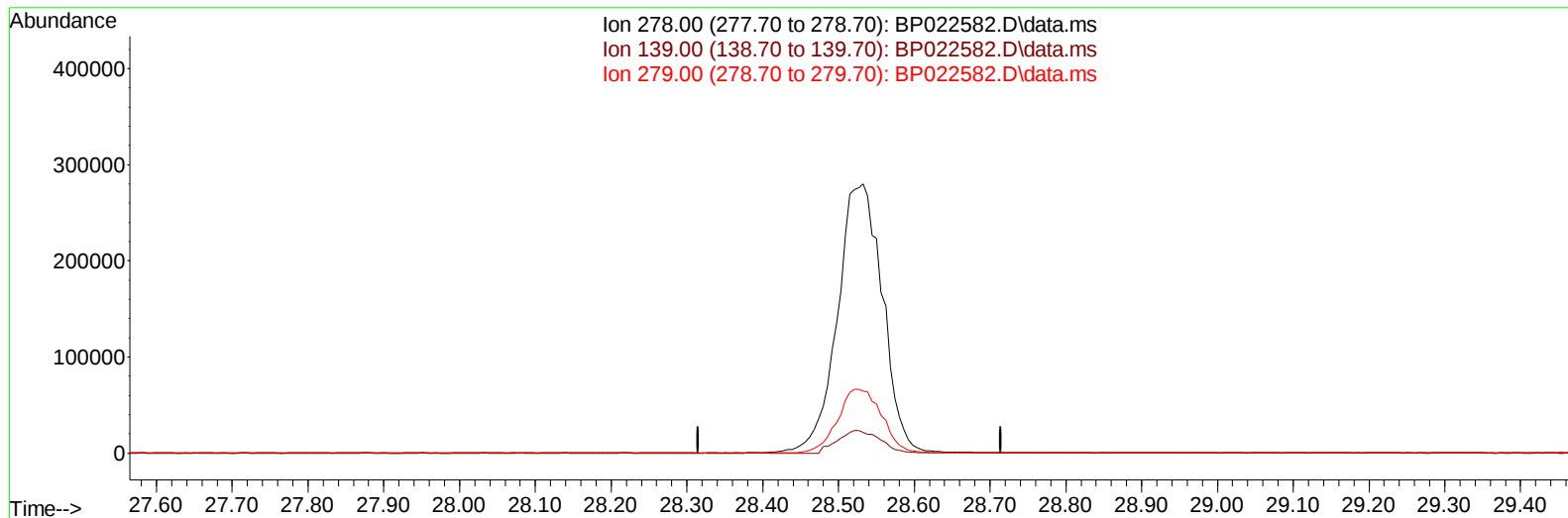
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Data File : BP022582.D  
Acq On : 24 Oct 2024 11:00  
Operator : RC/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:23:50 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 10/25/2024  
Supervised By :mohammad ahmed 10/26/2024



TIC: BP022582.D\data.ms

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

**28.515min (-28.515) 0.00 ng/u1**

**response 0**

| Ion    | Exp%   | Act%  |
|--------|--------|-------|
| 278.00 | 100.00 | 0.00  |
| 139.00 | 9.10   | 0.00# |
| 279.00 | 23.70  | 0.00# |
| 0.00   | 0.00   | 0.00  |

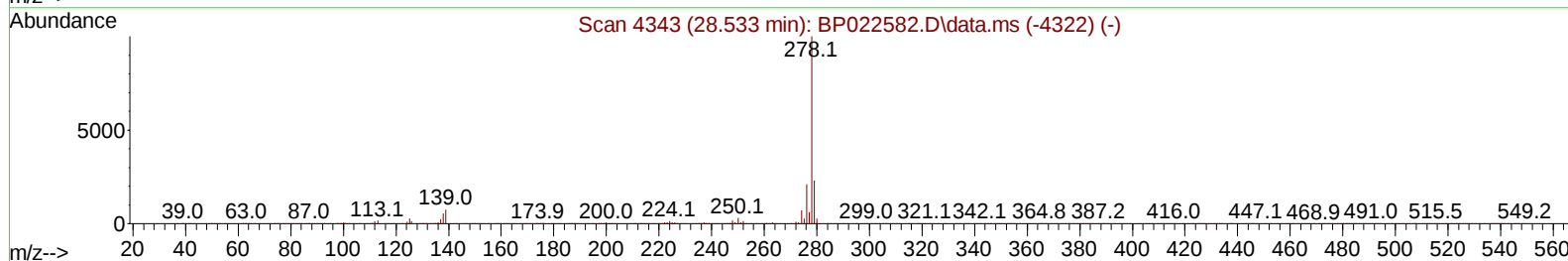
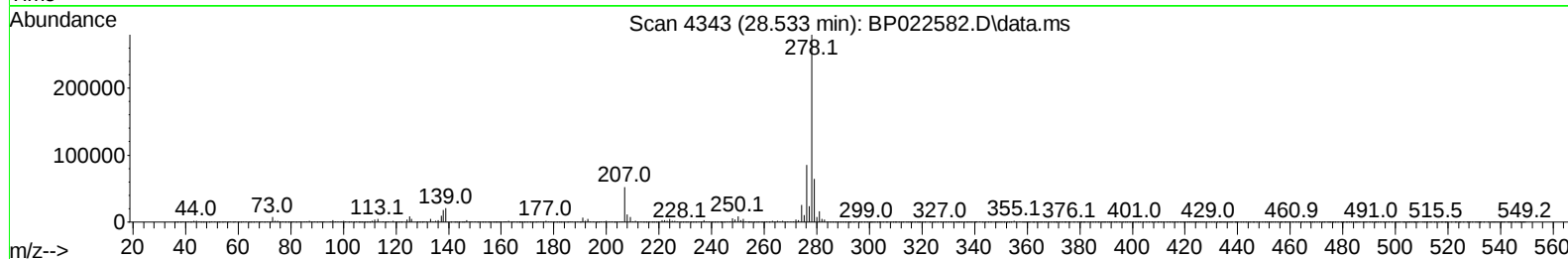
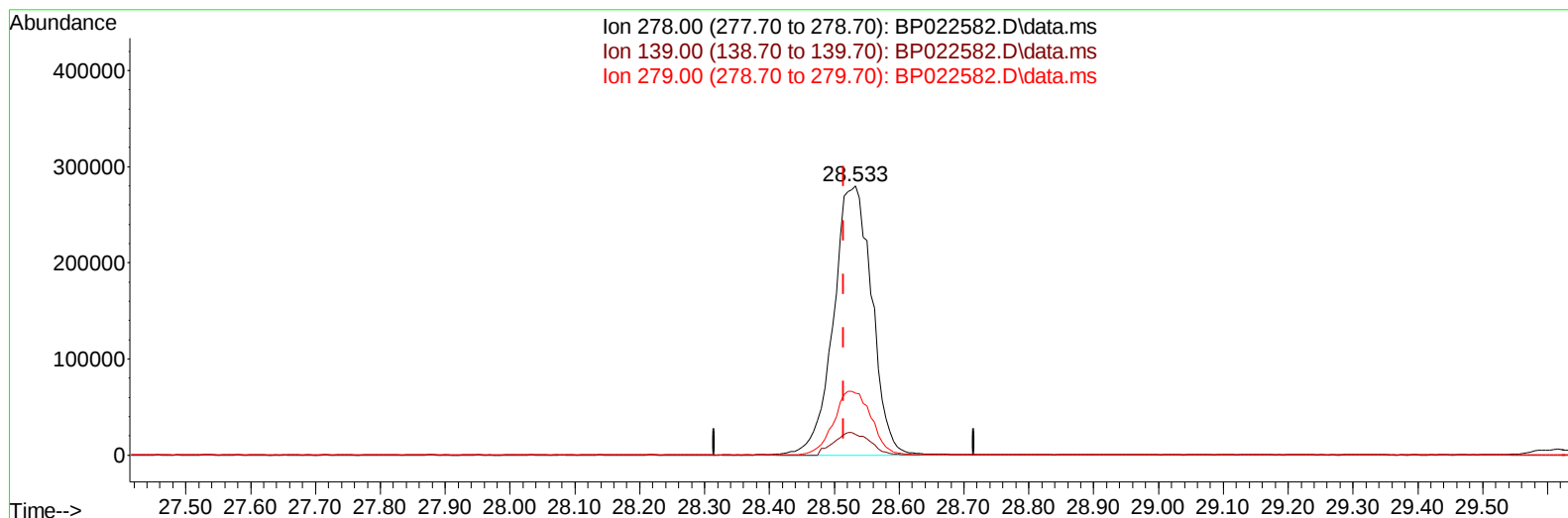
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Data File : BP022582.D  
Acq On : 24 Oct 2024 11:00  
Operator : RC/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:23:50 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 10/25/2024  
Supervised By :mohammad ahmed 10/26/2024



TIC: BP022582.D\data.ms

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

**28.533min (+ 0.018) 19.82 ng/ul m**

**response 1149960**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 278.00 | 100.00 | 100.00 |
| 139.00 | 9.10   | 7.60   |
| 279.00 | 23.70  | 22.92  |
| 0.00   | 0.00   | 0.00   |

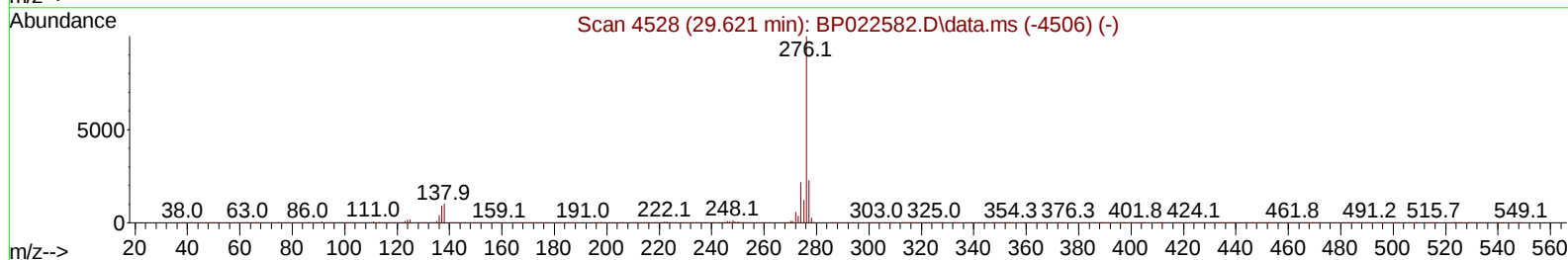
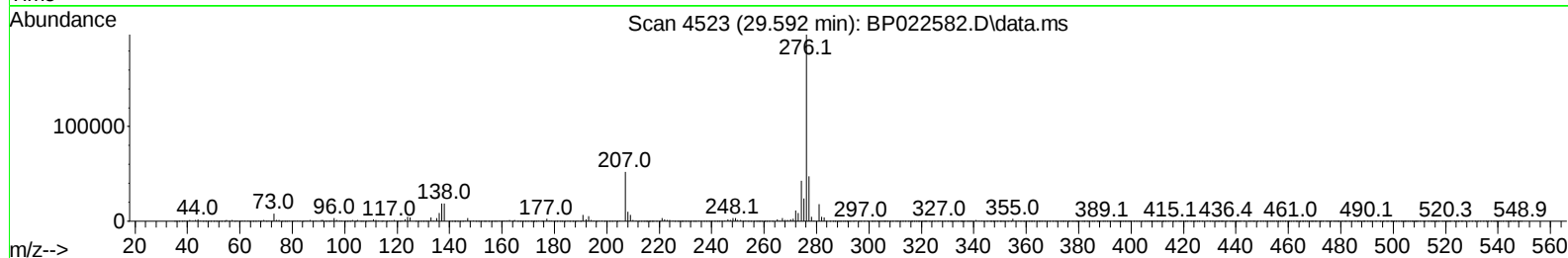
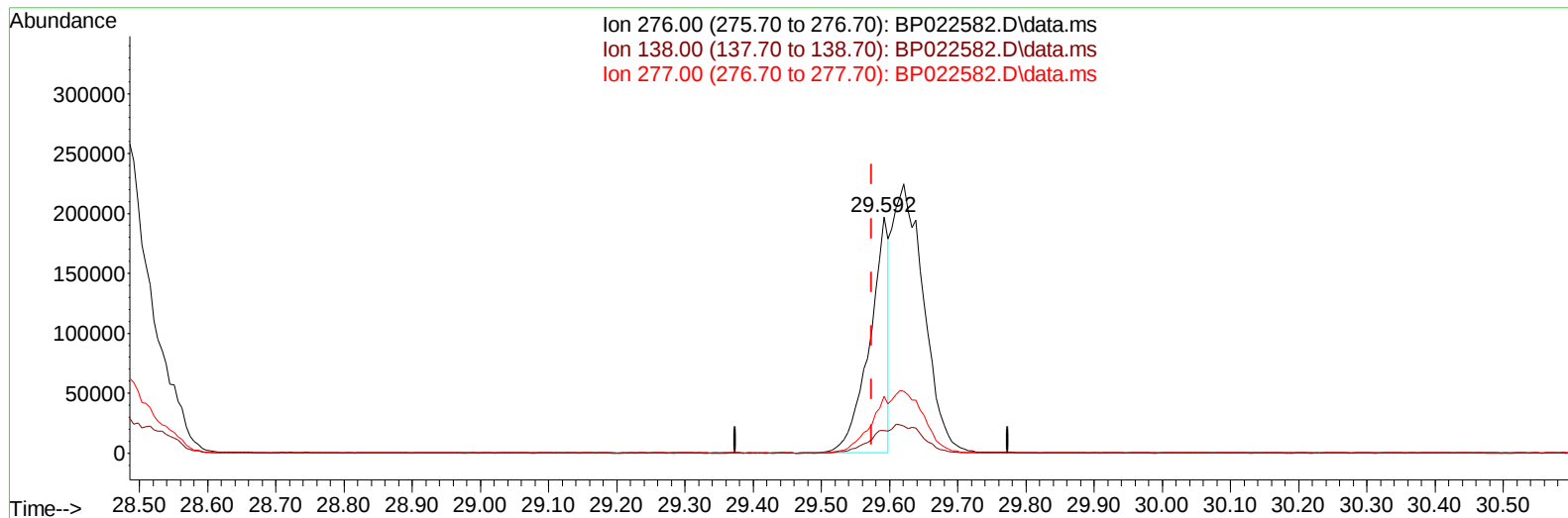
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102424\  
Data File : BP022582.D  
Acq On : 24 Oct 2024 11:00  
Operator : RC/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:23:50 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 10/25/2024  
Supervised By :mohammad ahmed 10/26/2024



TIC: BP022582.D\data.ms

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

**29.592min (+ 0.018) 6.87 ng/u1**

**response 383036**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 10.60  | 9.61   |
| 277.00 | 23.70  | 24.01  |
| 0.00   | 0.00   | 0.00   |

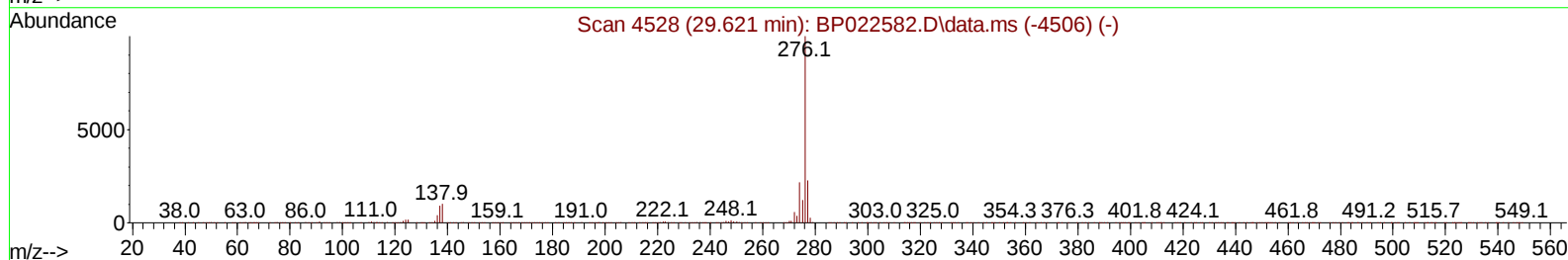
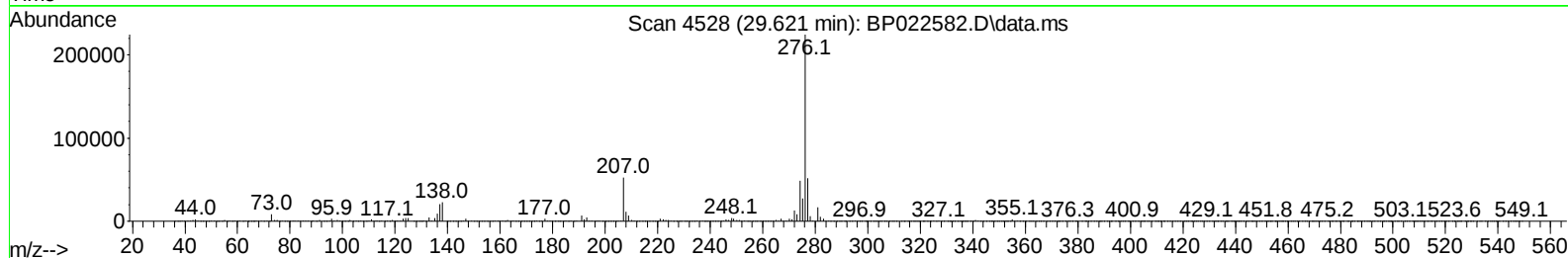
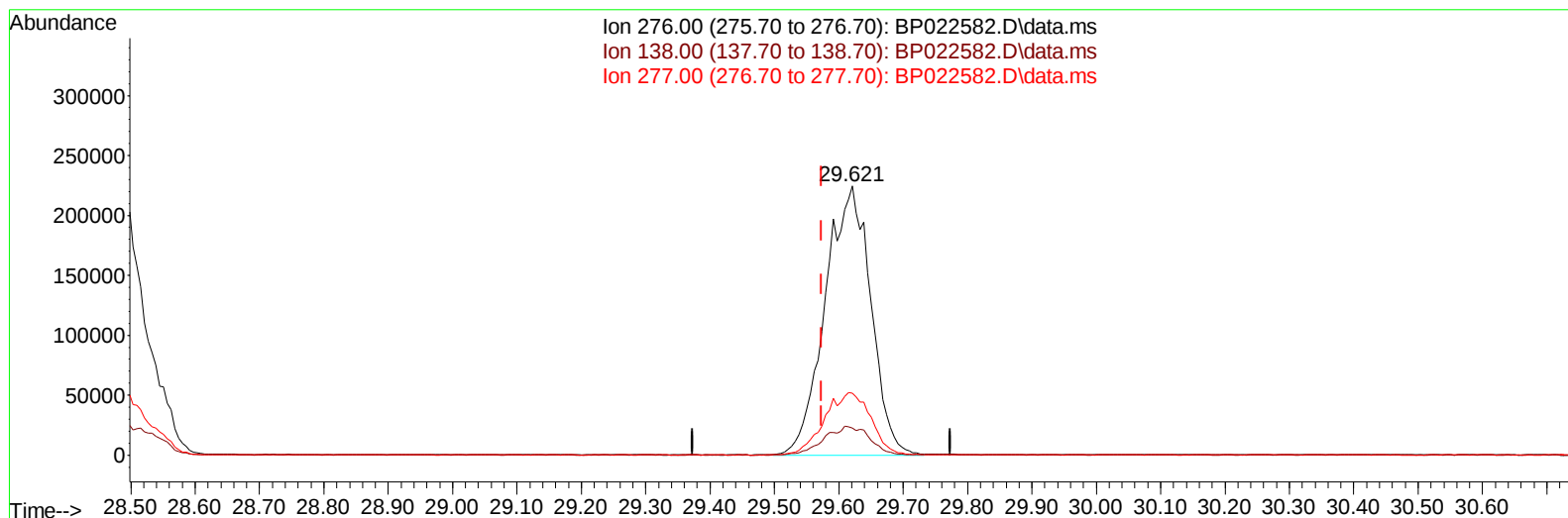
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102424\  
Data File : BP022582.D  
Acq On : 24 Oct 2024 11:00  
Operator : RC/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:23:50 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 10/25/2024  
Supervised By :mohammad ahmed 10/26/2024



TIC: BP022582.D\data.ms

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

29.621min (+ 0.047) 19.65 ng/ul m

response 1095424

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 10.60  | 10.08  |
| 277.00 | 23.70  | 22.90  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102424\  
 Data File : BP022582.D  
 Acq On : 24 Oct 2024 11:00  
 Operator : RC/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:24:46 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 10/25/2024  
 Supervised By :mohammad ahmed 10/26/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.652  | 152  | 116396   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.404 | 136  | 449747   | 20.000 | ng/ul  | -0.01    |
| 38) Acenaphthene-d10          | 14.269 | 164  | 352793   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.063 | 188  | 843139   | 20.000 | ng/ul  | -0.02    |
| 79) Chrysene-d12              | 21.510 | 240  | 834279   | 20.000 | ng/ul  | 0.01     |
| 88) Perylene-d12              | 24.768 | 264  | 926977   | 20.000 | ng/ul  | 0.02     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.187  | 96   | 20461    | 6.831  | ng/uL  | -0.01    |
| 4) Pyridine-d5                | 3.605  | 84   | 156845   | 19.075 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.846  | 99   | 193247   | 19.723 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.993  | 67   | 115114   | 19.990 | ng/ul  | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.193  | 132  | 145827   | 20.975 | ng/ul  | -0.01    |
| 15) 4-Methylphenol-d8         | 8.352  | 113  | 161913   | 20.707 | ng/ul  | -0.01    |
| 21) Nitrobenzene-d5           | 8.793  | 128  | 74195    | 21.455 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.499  | 143  | 90096    | 22.504 | ng/ul  | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.040 | 165  | 183620   | 21.579 | ng/ul  | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.540 | 131  | 205345   | 20.423 | ng/ul  | -0.02    |
| 46) Dimethylphthalate-d6      | 13.681 | 166  | 584675   | 20.859 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.957 | 160  | 625939   | 21.025 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.498 | 143  | 91180    | 19.390 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.269 | 176  | 494121   | 20.324 | ng/ul  | -0.02    |
| 65) 4,6-Dinitro-2-methylph... | 15.422 | 200  | 110261   | 19.884 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.175 | 188  | 821257   | 20.706 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.551 | 212  | 985961   | 22.348 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.545 | 264  | 1040381  | 21.016 | ng/ul  | 0.02     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.222  | 88   | 22402    | 6.956  | ng/uL  | 97       |
| 5) Pyridine                   | 3.622  | 79   | 159408   | 18.907 | ng/ul  | 99       |
| 6) Benzaldehyde               | 6.810  | 77   | 113841   | 21.501 | ng/ul  | 97       |
| 8) Phenol                     | 6.869  | 94   | 199885   | 19.608 | ng/ul  | 100      |
| 10) Bis(2-Chloroethyl)ether   | 7.087  | 93   | 154545   | 20.257 | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.228  | 128  | 151791   | 21.114 | ng/ul  | 97       |
| 13) 2-Methylphenol            | 8.093  | 108  | 152881   | 20.666 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.157  | 45   | 179994   | 20.300 | ng/ul  | 100      |
| 16) Acetophenone              | 8.457  | 105  | 266994   | 20.683 | ng/ul  | 96       |
| 17) N-Nitroso-di-n-propyla... | 8.446  | 70   | 139267   | 21.294 | ng/ul  | 97       |
| 18) 4-Methylphenol            | 8.416  | 108  | 164914   | 20.094 | ng/ul  | 96       |
| 19) Hexachloroethane          | 8.710  | 117  | 68179    | 20.363 | ng/ul  | 99       |
| 22) Nitrobenzene              | 8.834  | 77   | 204731   | 19.834 | ng/ul  | 98       |
| 23) Isophorone                | 9.351  | 82   | 389381   | 21.039 | ng/ul  | 100      |
| 25) 2-Nitrophenol             | 9.534  | 139  | 94582    | 21.920 | ng/ul  | 96       |
| 26) 2,4-Dimethylphenol        | 9.599  | 107  | 201175   | 21.451 | ng/ul  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 9.822  | 93   | 222601   | 21.183 | ng/ul  | 100      |
| 29) 2,4-Dichlorophenol        | 10.063 | 162  | 182664   | 21.799 | ng/ul  | 99       |
| 30) Naphthalene               | 10.457 | 128  | 499782   | 20.864 | ng/ul  | 98       |
| 32) 4-Chloroaniline           | 10.569 | 127  | 203265   | 20.480 | ng/ul  | 96       |
| 33) Hexachlorobutadiene       | 10.734 | 225  | 146900   | 20.800 | ng/ul  | 99       |
| 34) Caprolactam               | 11.351 | 113  | 53713m   | 19.883 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.704 | 107  | 193417   | 21.156 | ng/ul  | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102424\  
 Data File : BP022582.D  
 Acq On : 24 Oct 2024 11:00  
 Operator : RC/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:24:46 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 10/25/2024  
 Supervised By :mohammad ahmed 10/26/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.063 | 142  | 371612   | 21.134 | ng/ul | 96       |
| 37) 1-Methylnaphthalene       | 12.287 | 142  | 374337   | 20.857 | ng/ul | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.440 | 216  | 273045   | 20.931 | ng/ul | 95       |
| 40) Hexachlorocyclopentadiene | 12.416 | 237  | 127509   | 16.043 | ng/ul | 98       |
| 41) 2,4,6-Trichlorophenol     | 12.692 | 196  | 175384   | 21.374 | ng/ul | 97       |
| 42) 2,4,5-Trichlorophenol     | 12.769 | 196  | 188770   | 21.606 | ng/ul | 97       |
| 43) 1,1'-Biphenyl             | 13.087 | 154  | 532878   | 20.955 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.128 | 162  | 422229   | 20.283 | ng/ul | 98       |
| 45) 2-Nitroaniline            | 13.345 | 65   | 126856   | 20.409 | ng/ul | 99       |
| 47) Dimethylphthalate         | 13.728 | 163  | 571933   | 20.477 | ng/ul | 99       |
| 48) 2,6-Dinitrotoluene        | 13.845 | 165  | 113540   | 21.491 | ng/ul | 96       |
| 50) Acenaphthylene            | 13.987 | 152  | 635882   | 20.580 | ng/ul | 99       |
| 51) 3-Nitroaniline            | 14.175 | 138  | 97083    | 19.565 | ng/ul | 99       |
| 52) Acenaphthene              | 14.334 | 153  | 449561   | 20.576 | ng/ul | 100      |
| 53) 2,4-Dinitrophenol         | 14.410 | 184  | 72163    | 18.565 | ng/ul | 99       |
| 55) 4-Nitrophenol             | 14.510 | 109  | 98463    | 18.866 | ng/ul | 97       |
| 56) Dibenzofuran              | 14.675 | 168  | 671735   | 20.467 | ng/ul | 98       |
| 57) 2,4-Dinitrotoluene        | 14.651 | 165  | 177546   | 21.627 | ng/ul | 97       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.904 | 232  | 177589   | 21.777 | ng/ul | 98       |
| 59) Diethylphthalate          | 15.110 | 149  | 571868   | 21.341 | ng/ul | 99       |
| 61) Fluorene                  | 15.328 | 166  | 549647   | 20.554 | ng/ul | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.322 | 204  | 318469   | 20.825 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.357 | 138  | 99127    | 19.853 | ng/ul | 99       |
| 66) 4,6-Dinitro-2-methylph... | 15.434 | 198  | 118017   | 19.759 | ng/ul | 95       |
| 67) N-Nitrosodiphenylamine    | 15.539 | 169  | 475259   | 21.049 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.239 | 248  | 221059   | 21.742 | ng/ul | 97       |
| 69) Hexachlorobenzene         | 16.351 | 284  | 267579   | 21.380 | ng/ul | 99       |
| 70) Atrazine                  | 16.528 | 200  | 205725   | 21.920 | ng/ul | 94       |
| 71) Pentachlorophenol         | 16.716 | 266  | 165654   | 20.585 | ng/ul | 96       |
| 72) Phenanthrene              | 17.110 | 178  | 912011   | 20.218 | ng/ul | 100      |
| 74) Anthracene                | 17.210 | 178  | 925463   | 20.559 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.057 | 216  | 281933   | 21.961 | ng/uL | 98       |
| 76) Pentachlorobenzene        | 14.592 | 250  | 302327   | 21.303 | ng/uL | 99       |
| 77) Carbazole                 | 17.492 | 167  | 806885   | 20.257 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 18.069 | 149  | 980179   | 22.240 | ng/ul | 100      |
| 80) Fluoranthene              | 19.198 | 202  | 1166103  | 22.991 | ng/ul | 100      |
| 82) Pyrene                    | 19.580 | 202  | 1209913  | 22.257 | ng/ul | 99       |
| 83) Butylbenzylphthalate      | 20.533 | 149  | 441187   | 23.311 | ng/ul | 96       |
| 84) 3,3'-Dichlorobenzidine    | 21.404 | 252  | 416250   | 20.930 | ng/ul | 98       |
| 85) Benzo(a)anthracene        | 21.492 | 228  | 1191351  | 20.370 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.416 | 149  | 638626   | 23.508 | ng/ul | 99       |
| 87) Chrysene                  | 21.557 | 228  | 1104211  | 20.362 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.651 | 149  | 1055181  | 23.931 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.733 | 252  | 1208264  | 20.707 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.804 | 252  | 1211541  | 21.208 | ng/ul | 99       |
| 93) Benzo(a)pyrene            | 24.610 | 252  | 1100660  | 20.496 | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.468 | 276  | 1431795  | 19.979 | ng/ul | 99       |
| 95) Dibenzo(a,h)anthracene    | 28.533 | 278  | 1149960m | 19.816 | ng/ul |          |
| 96) Benzo(g,h,i)perylene      | 29.621 | 276  | 1095424m | 19.652 | ng/ul |          |

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

**Instrument :**  
BNA\_P  
**LabSampleId :**  
SSTDCCC020

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 10/25/2024  
Supervised By :mohammad ahmed 10/26/2024

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102424\  
 Data File : BP022582.D  
 Acq On : 24 Oct 2024 11:00  
 Operator : RC/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 24 12:24:46 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 10/25/2024  
 Supervised By :mohammad ahmed 10/26/2024

