

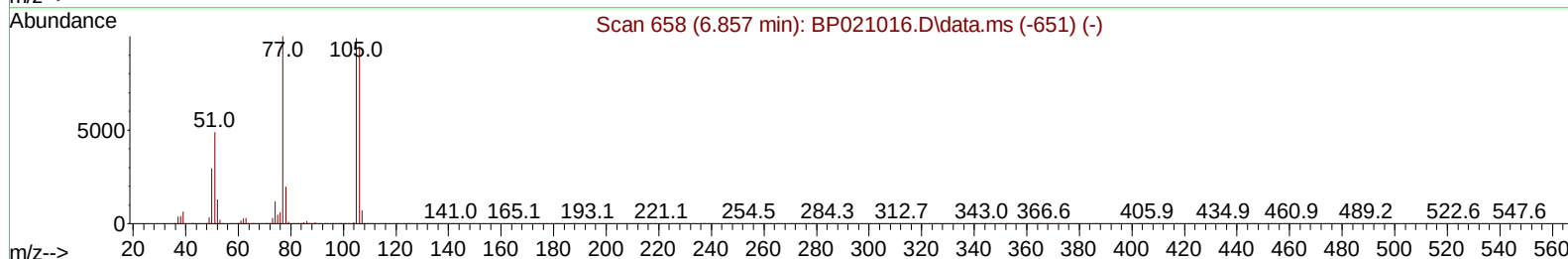
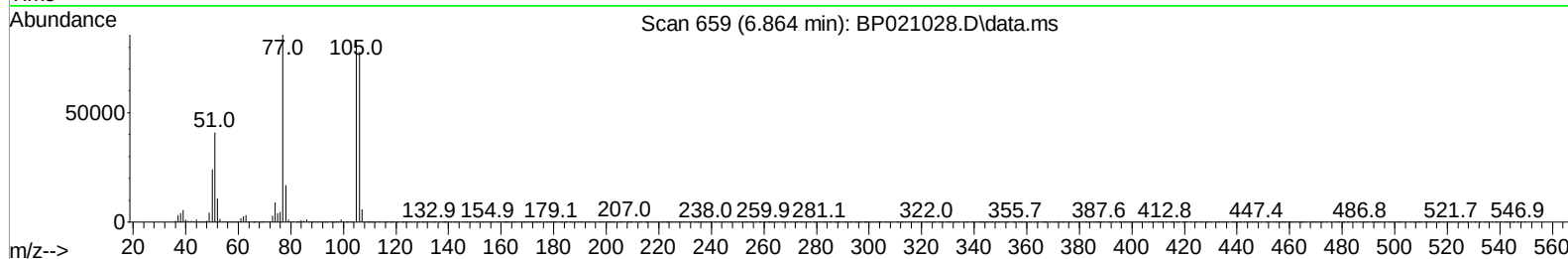
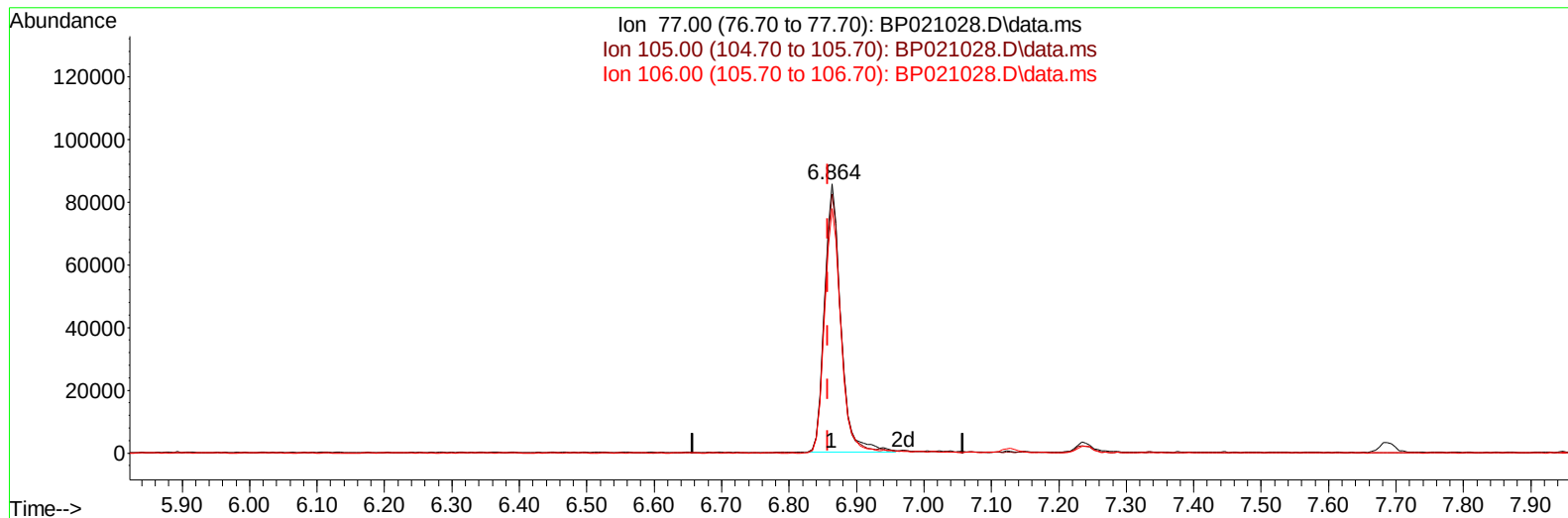
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP071724\  
 Data File : BP021028.D  
 Acq On : 17 Jul 2024 23:15  
 Operator : MA/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Jul 18 00:40:52 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 15 12:22:22 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/18/2024  
 Supervised By :mohammad ahmed 07/19/2024



TIC: BP021028.D\data.ms

**(6) Benzaldehyde**

**6.864min (+ 0.006) 23.32 ng/u1**

**response 144478**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 99.10  | 96.22  |
| 106.00 | 93.30  | 90.81  |
| 0.00   | 0.00   | 0.00   |

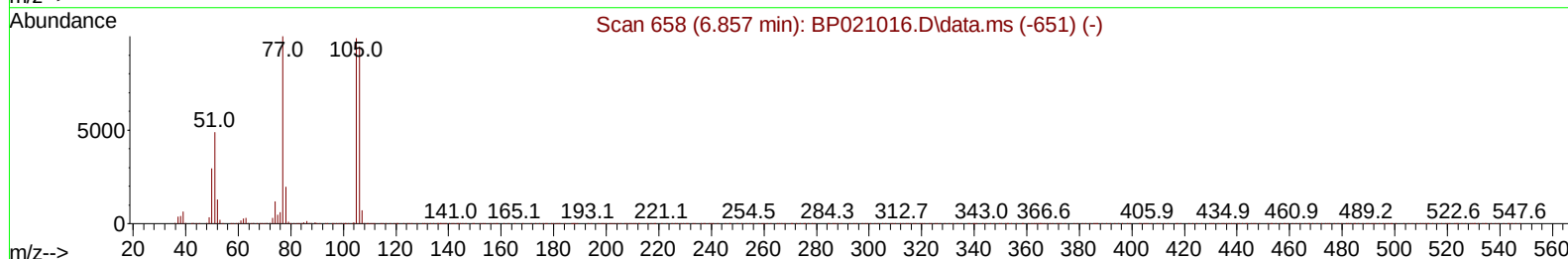
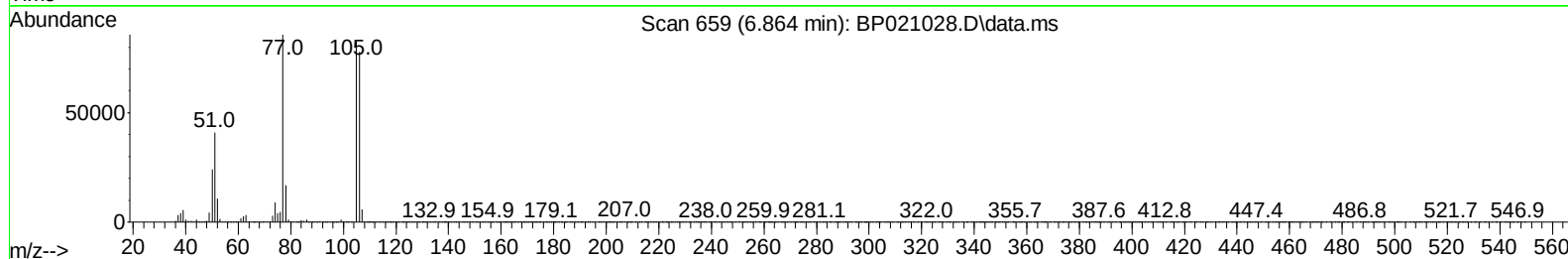
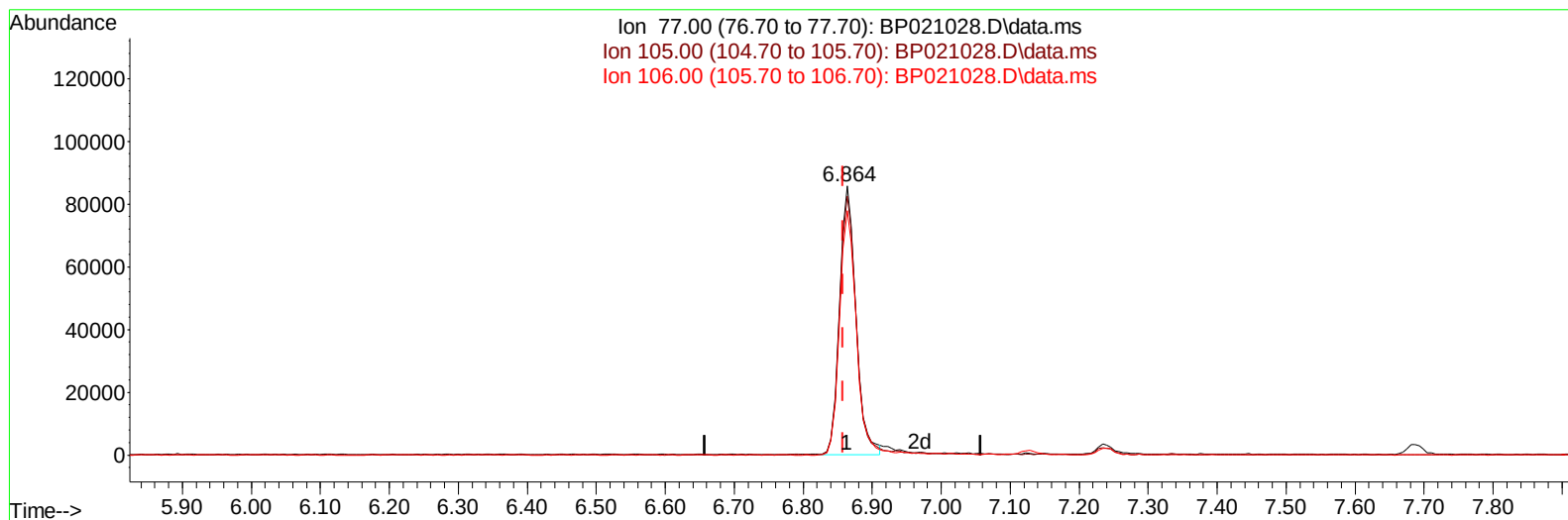
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TIC: BP021028.D\data.ms

(6) Benzaldehyde

6.864min (+ 0.006) 22.78 ng/ul m

response 141174

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 99.10  | 96.22  |
| 106.00 | 93.30  | 90.81  |
| 0.00   | 0.00   | 0.00   |

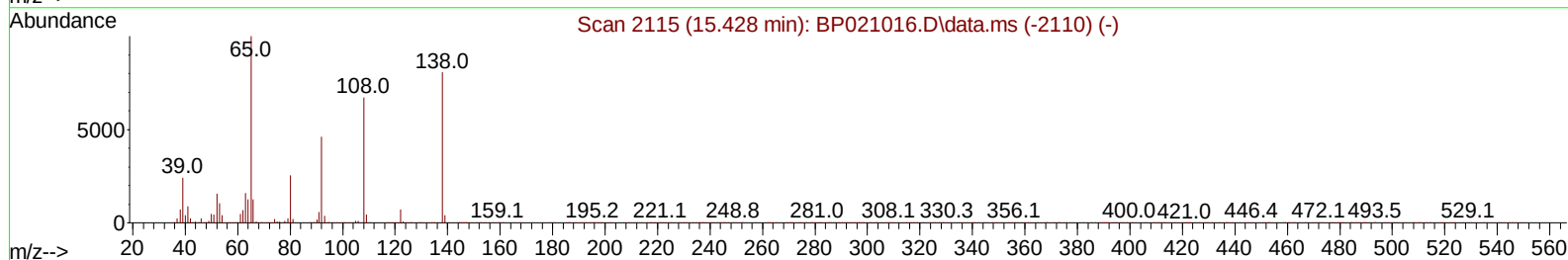
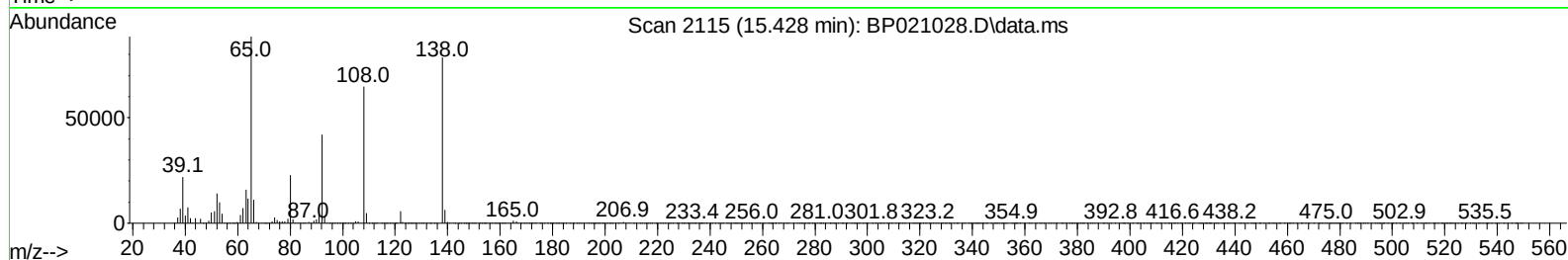
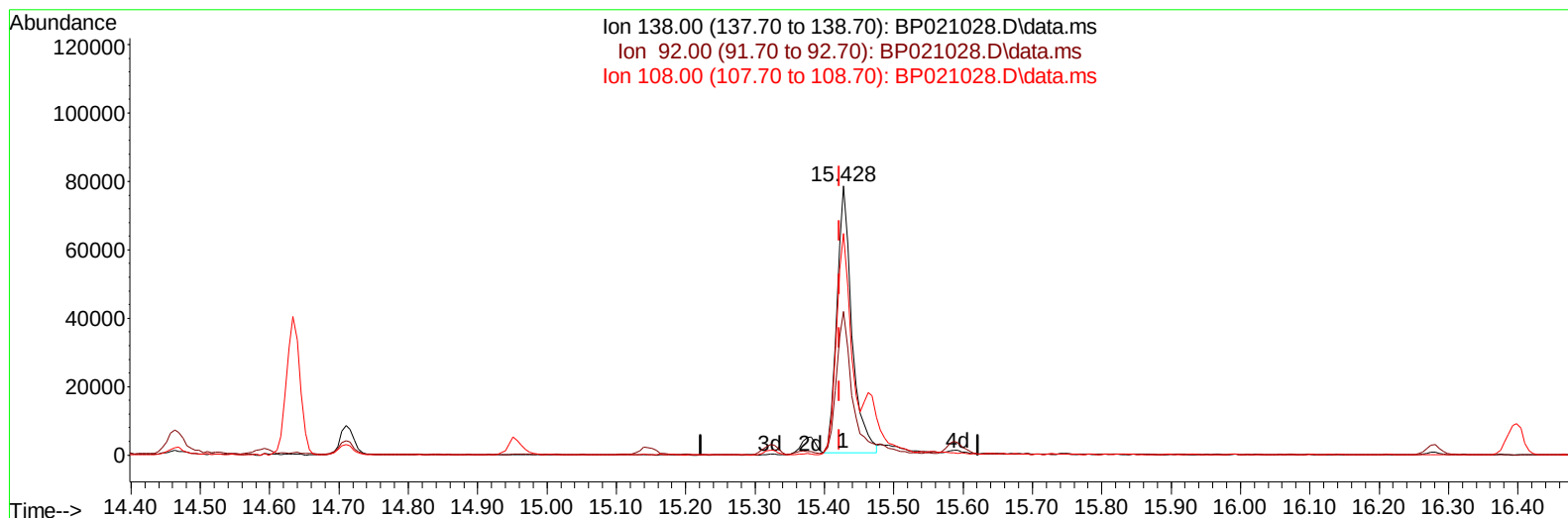
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TIC: BP021028.D\data.ms

**(63) 4-Nitroaniline**

**15.428min (+ 0.006) 26.46 ng/ul**

**response 116025**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 51.10  | 53.51  |
| 108.00 | 80.90  | 82.30  |
| 0.00   | 0.00   | 0.00   |

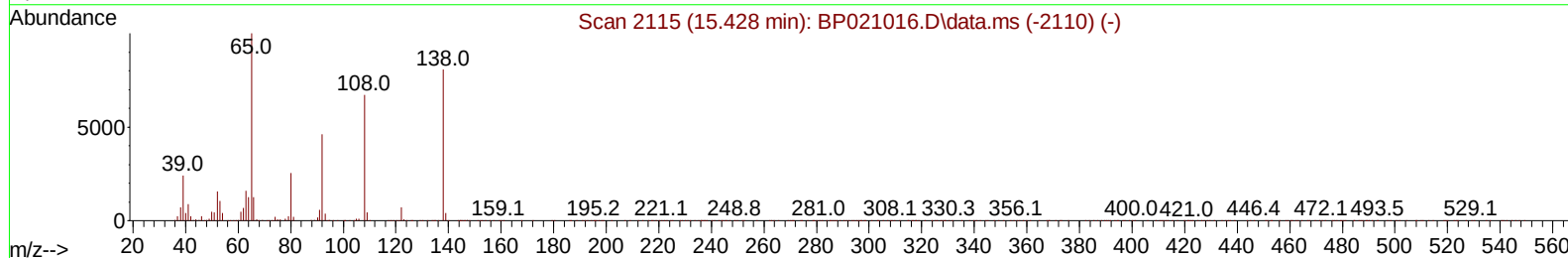
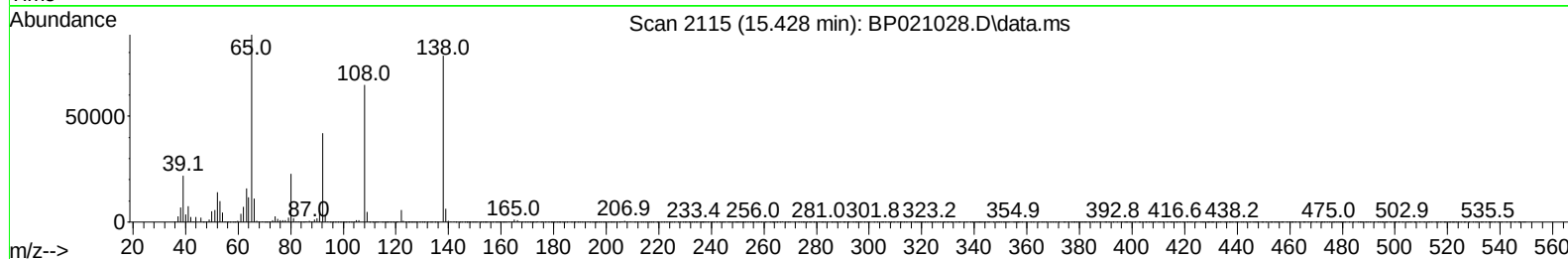
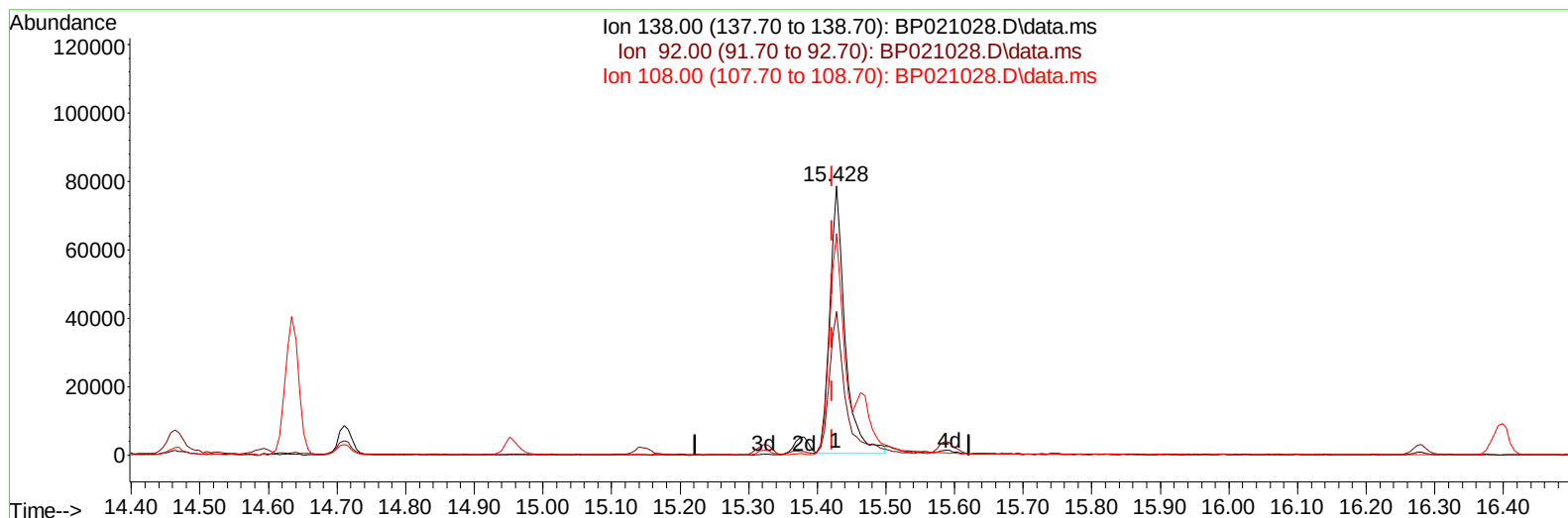
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Sample : SSTDCCC020  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

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Supervised By :mohammad ahmed 07/19/2024



TIC: BP021028.D\data.ms

(63) 4-Nitroaniline

15.428min (+ 0.006) 27.47 ng/ul m

response 120448

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 51.10  | 53.51  |
| 108.00 | 80.90  | 82.30  |
| 0.00   | 0.00   | 0.00   |

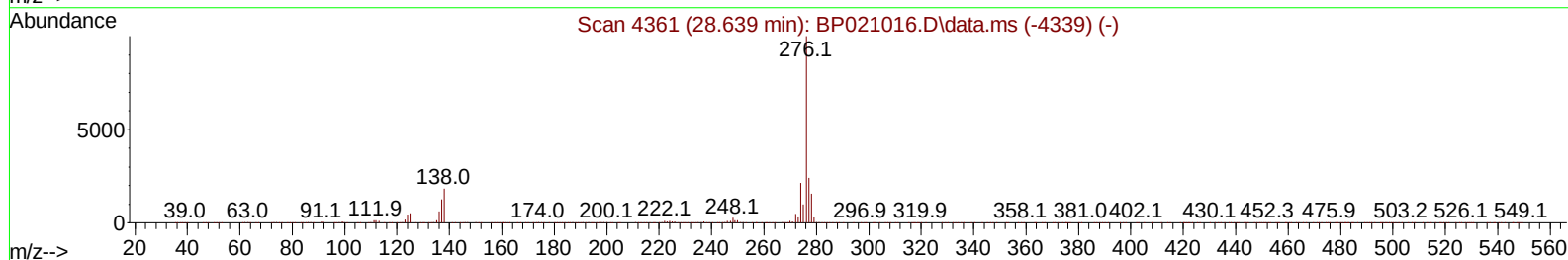
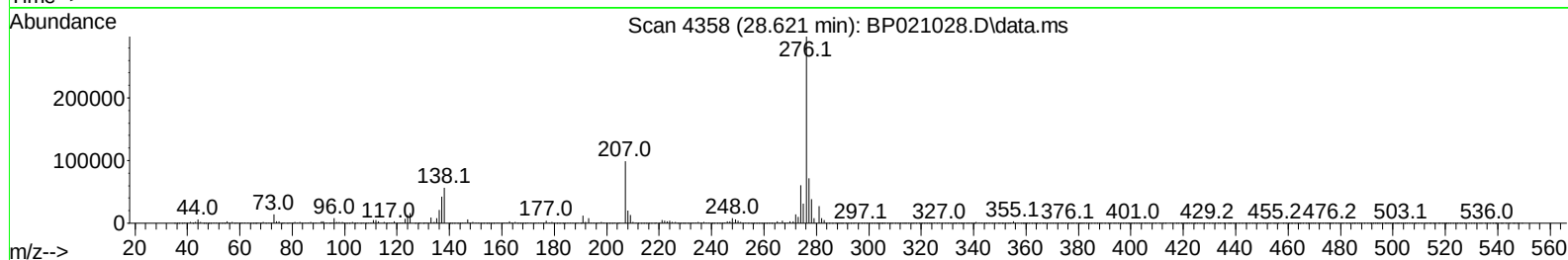
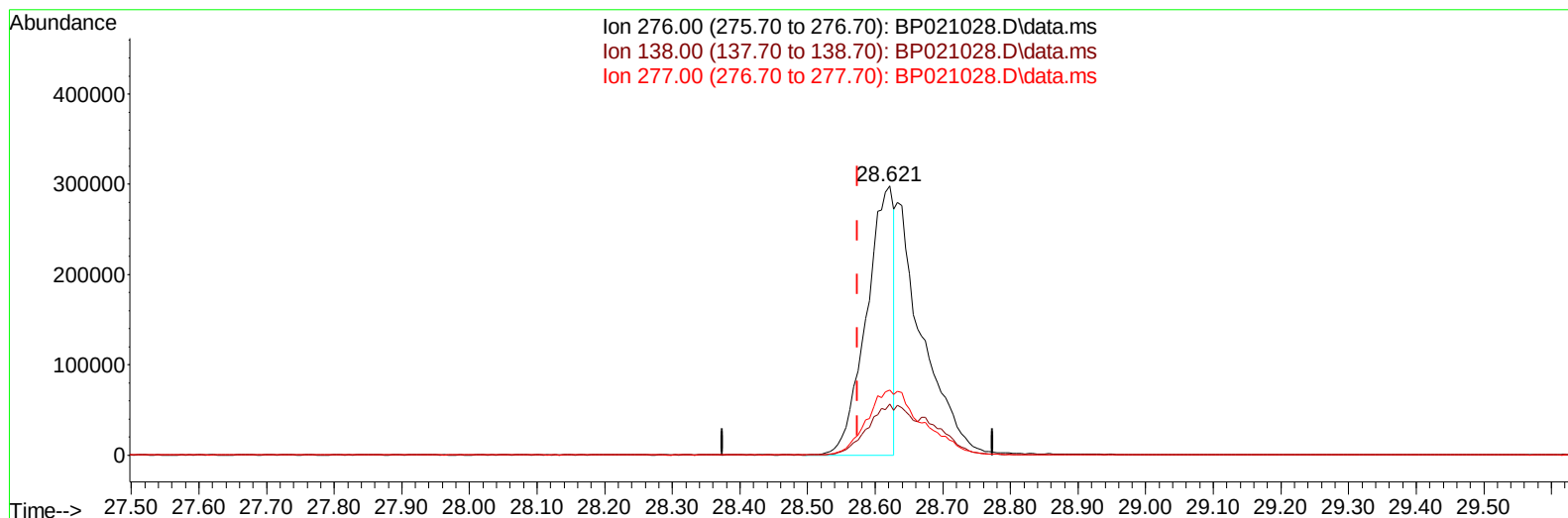
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Data File : BP021028.D  
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Sample : SSTDCCC020  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

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Reviewed By :Yogesh Patel 07/18/2024  
Supervised By :mohammad ahmed 07/19/2024



TIC: BP021028.D\data.ms

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

**28.621min (+ 0.047) 9.45 ng/u1**

**response 836170**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 17.90  | 19.09  |
| 277.00 | 24.00  | 24.12  |
| 0.00   | 0.00   | 0.00   |

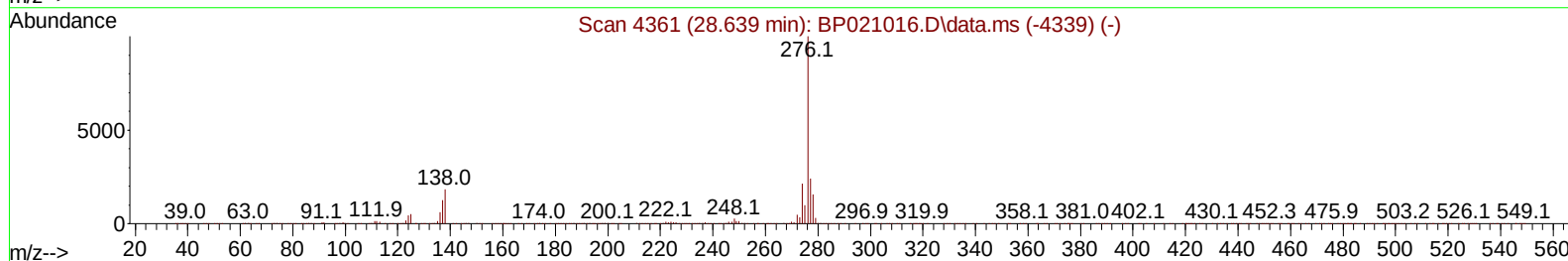
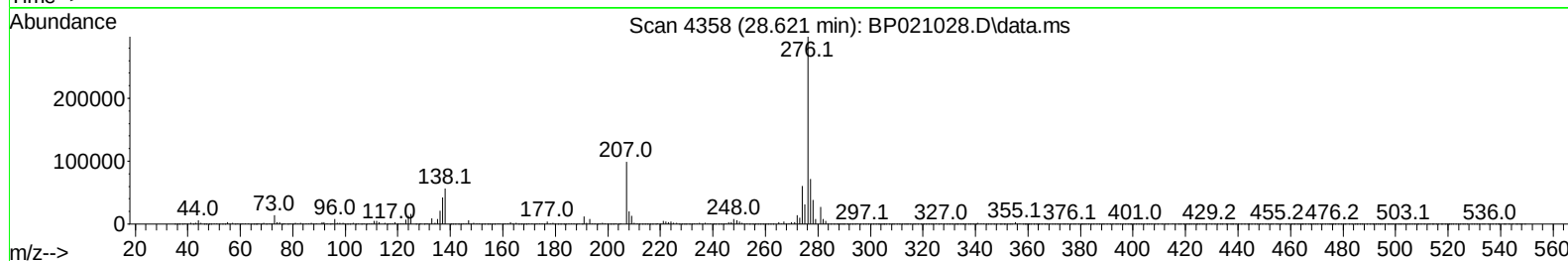
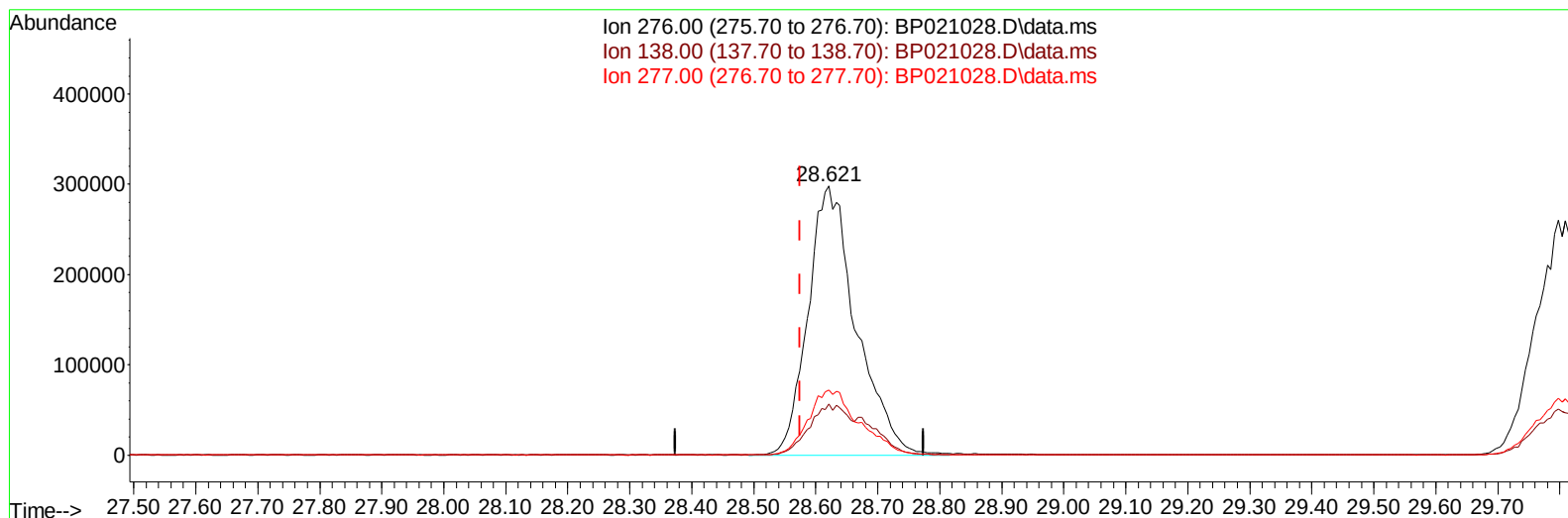
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP071724\  
Data File : BP021028.D  
Acq On : 17 Jul 2024 23:15  
Operator : MA/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Jul 18 00:40:52 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 15 12:22:22 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/18/2024  
Supervised By :mohammad ahmed 07/19/2024



TIC: BP021028.D\data.ms

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

**28.621min (+ 0.047) 18.19 ng/ul m**

**response 1608484**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 17.90  | 19.09  |
| 277.00 | 24.00  | 24.12  |
| 0.00   | 0.00   | 0.00   |

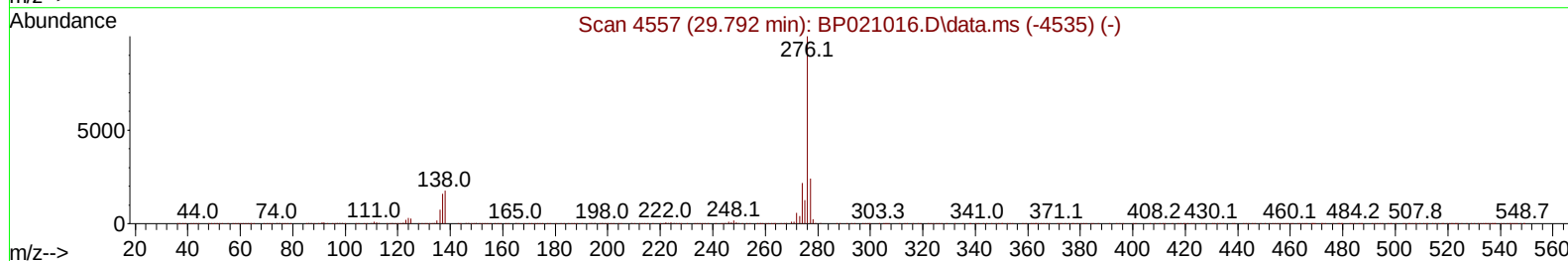
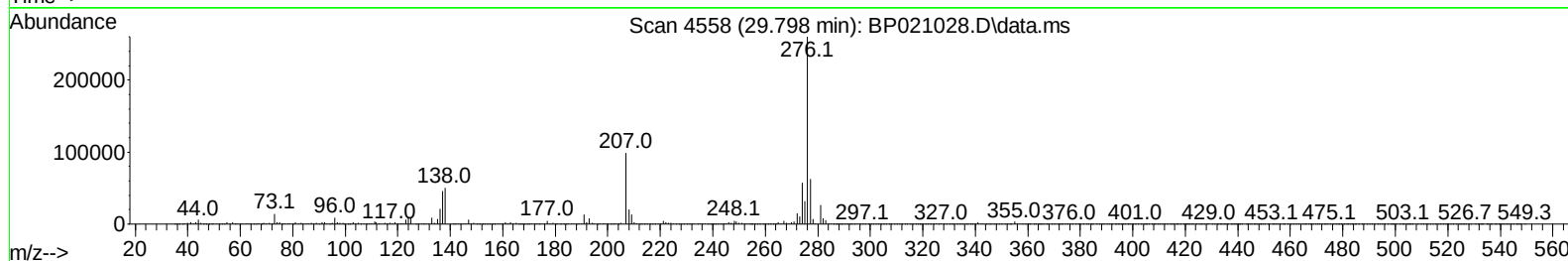
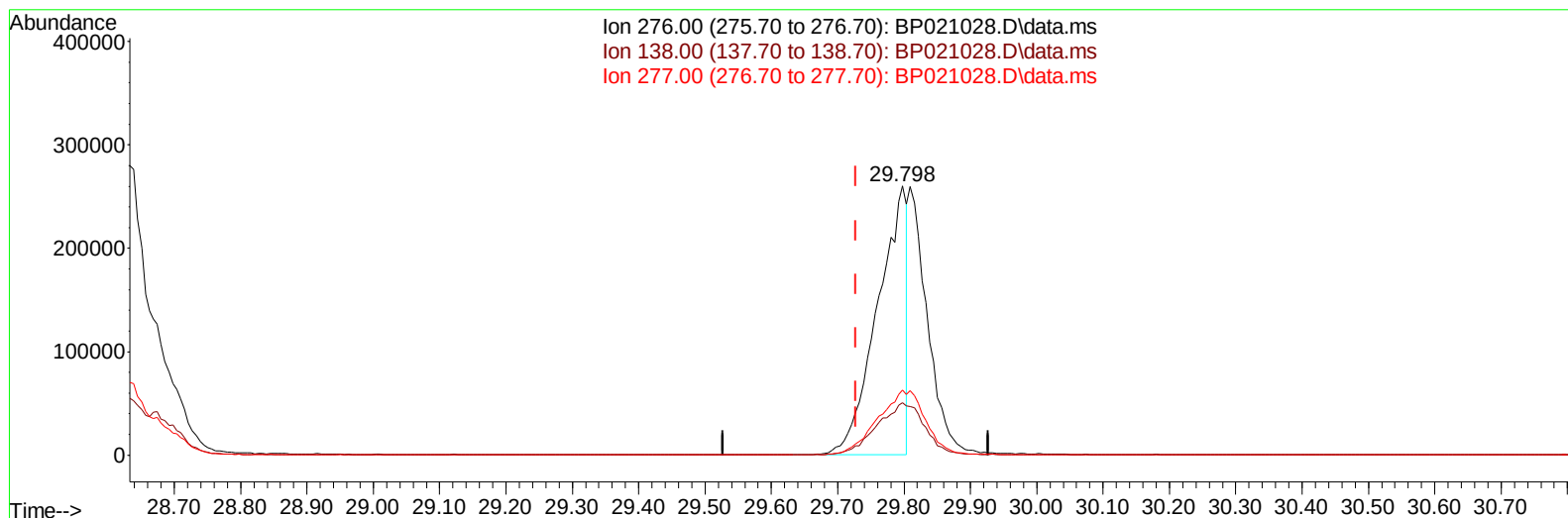
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Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

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TIC: BP021028.D\data.ms

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

29.798min (+ 0.071) 11.10 ng/ul

response 799382

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 19.10  | 19.48  |
| 277.00 | 22.70  | 24.17  |
| 0.00   | 0.00   | 0.00   |

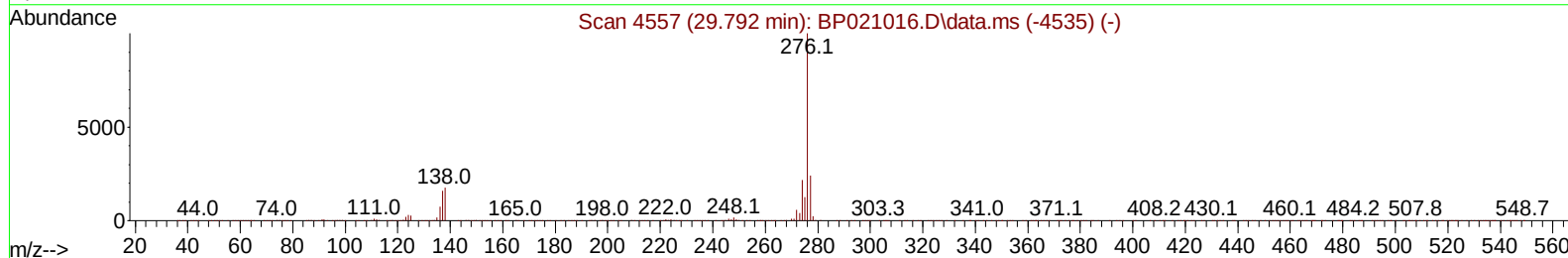
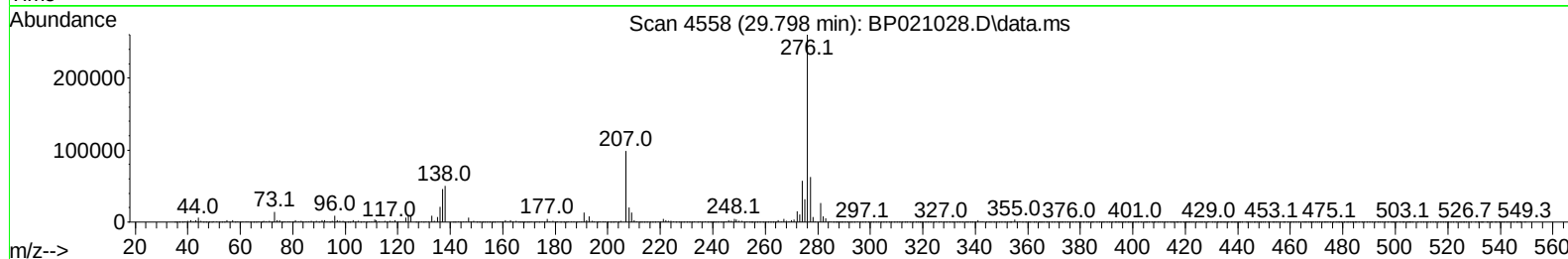
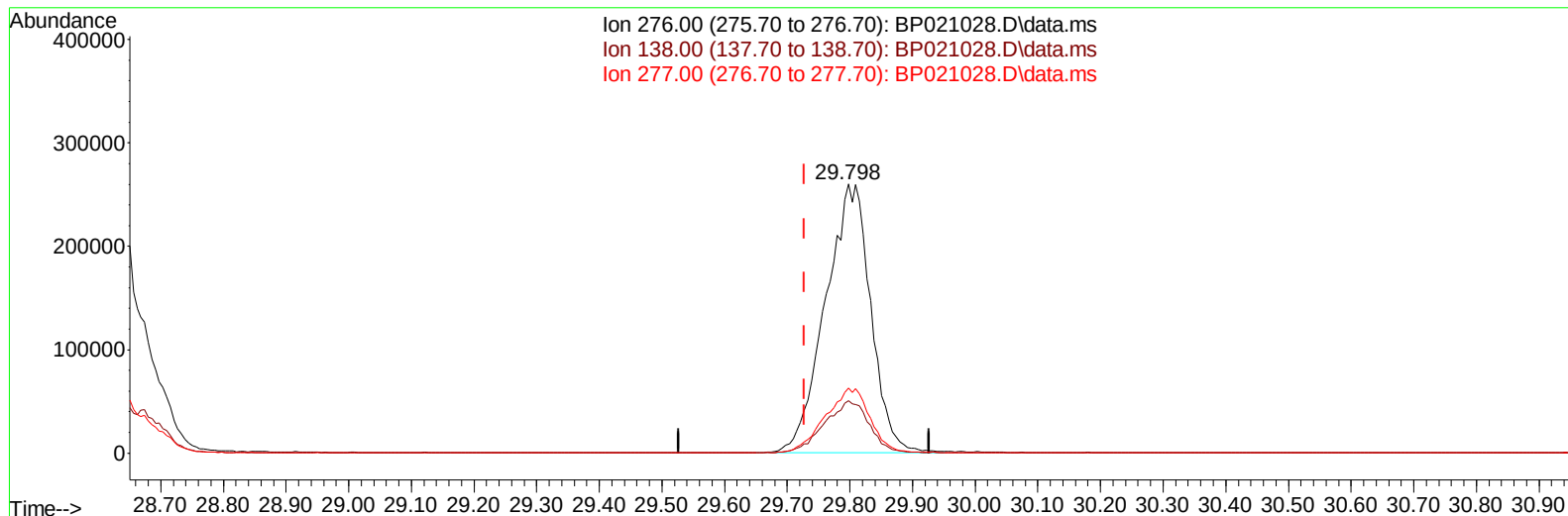
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Sample : SSTDCCC020  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020

Manual IntegrationsAPPROVED

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Reviewed By :Yogesh Patel 07/18/2024  
Supervised By :mohammad ahmed 07/19/2024



TIC: BP021028.D\data.ms

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

29.798min (+ 0.071) 18.20 ng/ul m

response 1310838

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 19.10  | 19.48  |
| 277.00 | 22.70  | 24.17  |
| 0.00   | 0.00   | 0.00   |



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Manual IntegrationsAPPROVED

Quant Time: Jul 18 01:05:12 2024  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.687  | 152  | 150762   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.452 | 136  | 624975   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.316 | 164  | 412604   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.122 | 188  | 949657   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.563 | 240  | 1015623  | 20.000 | ng/ul  | 0.01     |
| 88) Perylene-d12              | 24.880 | 264  | 1196822  | 20.000 | ng/ul  | 0.05     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.240  | 96   | 26724    | 8.066  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.658  | 84   | 190040   | 19.813 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.893  | 99   | 232595   | 18.924 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.034  | 67   | 135188   | 18.149 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.234  | 132  | 191147   | 18.875 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.399  | 113  | 186822   | 18.301 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.834  | 128  | 91353    | 18.819 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.546  | 143  | 102511   | 18.451 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.081 | 165  | 192898   | 19.201 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.599 | 131  | 261179   | 19.624 | ng/ul  | 0.01     |
| 46) Dimethylphthalate-d6      | 13.716 | 166  | 578141   | 19.275 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.998 | 160  | 655119   | 19.452 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.575 | 143  | 84394    | 19.776 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.322 | 176  | 498280   | 20.250 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.463 | 200  | 98041    | 17.063 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.228 | 188  | 829247   | 19.660 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.604 | 212  | 1029799  | 16.401 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.639 | 264  | 1120857  | 18.989 | ng/ul  | 0.04     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.276  | 88   | 28160    | 7.728  | ng/uL  | 96       |
| 5) Pyridine                   | 3.676  | 79   | 195497   | 19.761 | ng/ul  | 99       |
| 6) Benzaldehyde               | 6.864  | 77   | 141174m  | 22.783 | ng/ul  |          |
| 8) Phenol                     | 6.917  | 94   | 235091   | 18.902 | ng/ul  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.128  | 93   | 186643   | 18.166 | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.269  | 128  | 191918   | 18.909 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.140  | 108  | 179099   | 18.620 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.199  | 45   | 272640   | 18.367 | ng/ul  | 97       |
| 16) Acetophenone              | 8.505  | 105  | 289889   | 18.399 | ng/ul  | 97       |
| 17) N-Nitroso-di-n-propyla... | 8.475  | 70   | 147243   | 17.836 | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.463  | 108  | 192354   | 18.666 | ng/ul  | 96       |
| 19) Hexachloroethane          | 8.740  | 117  | 80681    | 17.689 | ng/ul  | 97       |
| 22) Nitrobenzene              | 8.881  | 77   | 220410   | 18.955 | ng/ul  | 97       |
| 23) Isophorone                | 9.399  | 82   | 423411   | 18.316 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.575  | 139  | 108863   | 18.787 | ng/ul  | 98       |
| 26) 2,4-Dimethylphenol        | 9.646  | 107  | 216300   | 18.930 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 9.869  | 93   | 255337   | 18.171 | ng/ul  | 98       |
| 29) 2,4-Dichlorophenol        | 10.110 | 162  | 188025   | 18.990 | ng/ul  | 99       |
| 30) Naphthalene               | 10.499 | 128  | 625366   | 19.023 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.622 | 127  | 245754   | 19.337 | ng/ul  | 100      |
| 33) Hexachlorobutadiene       | 10.757 | 225  | 132986   | 17.913 | ng/ul  | 98       |
| 34) Caprolactam               | 11.434 | 113  | 62671    | 20.547 | ng/ul  | 96       |
| 35) 4-Chloro-3-methylphenol   | 11.763 | 107  | 208416   | 19.365 | ng/ul  | 99       |

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 Supervised By :mohammad ahmed 07/19/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.110 | 142  | 416727   | 18.467 | ng/ul  | 98       |
| 37) 1-Methylnaphthalene       | 12.328 | 142  | 441683   | 19.031 | ng/ul  | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.481 | 216  | 251920   | 18.800 | ng/ul  | 97       |
| 40) Hexachlorocyclopentadiene | 12.446 | 237  | 110739   | 15.798 | ng/ul  | 97       |
| 41) 2,4,6-Trichlorophenol     | 12.740 | 196  | 161960   | 19.128 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol     | 12.828 | 196  | 168729   | 18.835 | ng/ul  | 98       |
| 43) 1,1'-Biphenyl             | 13.134 | 154  | 565640   | 18.759 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.175 | 162  | 465657   | 19.344 | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.404 | 65   | 135567   | 20.365 | ng/ul  | 97       |
| 47) Dimethylphthalate         | 13.763 | 163  | 586928   | 19.285 | ng/ul  | 100      |
| 48) 2,6-Dinitrotoluene        | 13.898 | 165  | 118712   | 19.510 | ng/ul  | 98       |
| 50) Acenaphthylene            | 14.028 | 152  | 731164   | 19.174 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.240 | 138  | 120530   | 22.764 | ng/ul  | 98       |
| 52) Acenaphthene              | 14.375 | 153  | 493084   | 19.482 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.463 | 184  | 50611    | 15.839 | ng/ul  | 94       |
| 55) 4-Nitrophenol             | 14.593 | 109  | 67792    | 19.529 | ng/ul  | 93       |
| 56) Dibenzofuran              | 14.710 | 168  | 685625   | 19.782 | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene        | 14.710 | 165  | 177525   | 21.254 | ng/ul# | 82       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.951 | 232  | 151773   | 19.785 | ng/ul  | 96       |
| 59) Diethylphthalate          | 15.145 | 149  | 585732   | 19.258 | ng/ul  | 100      |
| 61) Fluorene                  | 15.375 | 166  | 564440   | 19.957 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.375 | 204  | 293082   | 19.097 | ng/ul  | 98       |
| 63) 4-Nitroaniline            | 15.428 | 138  | 120448m  | 27.469 | ng/ul  |          |
| 66) 4,6-Dinitro-2-methylph... | 15.481 | 198  | 105711   | 17.349 | ng/ul  | 99       |
| 67) N-Nitrosodiphenylamine    | 15.587 | 169  | 487784   | 17.639 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.281 | 248  | 188913   | 16.972 | ng/ul  | 98       |
| 69) Hexachlorobenzene         | 16.398 | 284  | 228554   | 17.908 | ng/ul  | 97       |
| 70) Atrazine                  | 16.569 | 200  | 189036   | 18.756 | ng/ul  | 99       |
| 71) Pentachlorophenol         | 16.775 | 266  | 125032   | 18.037 | ng/ul  | 98       |
| 72) Phenanthrene              | 17.169 | 178  | 952508   | 19.335 | ng/ul  | 99       |
| 74) Anthracene                | 17.257 | 178  | 954454   | 19.004 | ng/ul  | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.093 | 216  | 248539   | 16.352 | ng/uL  | 99       |
| 76) Pentachlorobenzene        | 14.634 | 250  | 253210   | 16.765 | ng/uL  | 99       |
| 77) Carbazole                 | 17.545 | 167  | 883166   | 21.511 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.116 | 149  | 1035279  | 18.277 | ng/ul  | 100      |
| 80) Fluoranthene              | 19.251 | 202  | 1204611  | 15.886 | ng/ul  | 98       |
| 82) Pyrene                    | 19.633 | 202  | 1265028  | 16.394 | ng/ul  | 100      |
| 83) Butylbenzylphthalate      | 20.580 | 149  | 525008   | 16.330 | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.469 | 252  | 422298   | 18.197 | ng/ul  | 99       |
| 85) Benzo(a)anthracene        | 21.539 | 228  | 1346689  | 18.929 | ng/ul  | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.457 | 149  | 750669   | 16.225 | ng/ul  | 99       |
| 87) Chrysene                  | 21.610 | 228  | 1233291  | 18.655 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.704 | 149  | 1349200  | 17.065 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.821 | 252  | 1340674  | 18.459 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.886 | 252  | 1367866  | 18.872 | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 24.704 | 252  | 1288562  | 18.775 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.621 | 276  | 1608484m | 18.187 | ng/ul  |          |
| 95) Dibenzo(a,h)anthracene    | 28.674 | 278  | 1333258  | 18.205 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 29.798 | 276  | 1310838m | 18.199 | ng/ul  |          |

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

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

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 07/18/2024  
Supervised By :mohammad ahmed 07/19/2024

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