

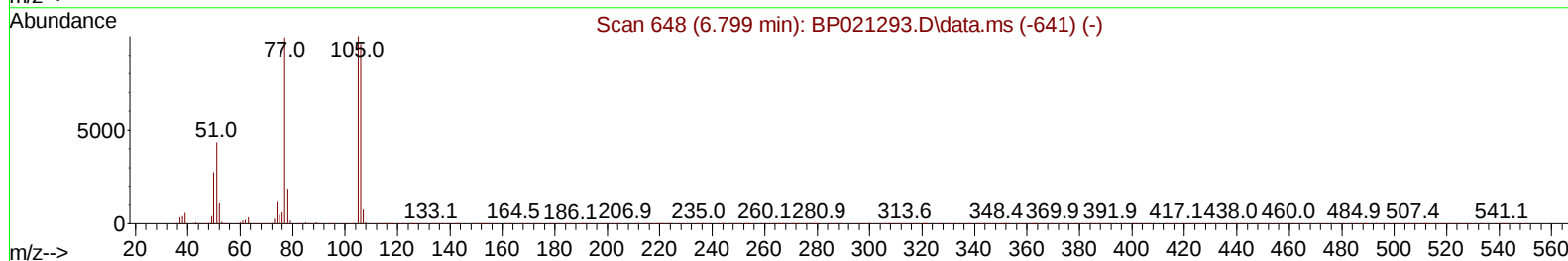
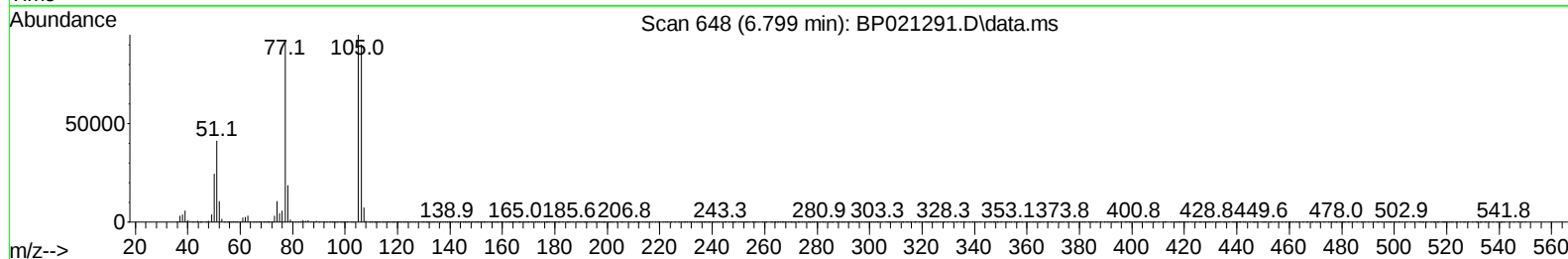
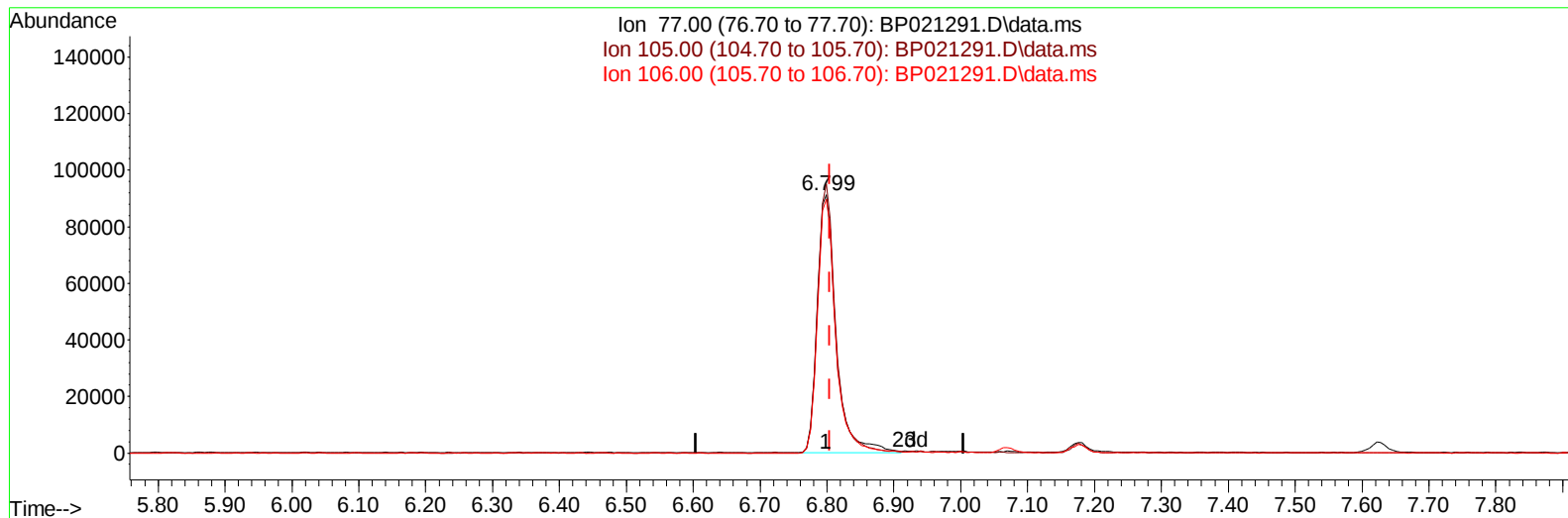
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072924\  
 Data File : BP021291.D  
 Acq On : 01 Aug 2024 13:24  
 Operator : CG/JU  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 LabSampleId :  
 SSTDCCC020EC

Manual IntegrationsAPPROVED

Quant Time: Aug 01 16:38:40 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 29 12:46:18 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/02/2024  
 Supervised By :mohammad ahmed 09/04/2024



TIC: BP021291.D\data.ms

**(6) Benzaldehyde**

**6.799min (-0.005) 21.11 ng/ul**

**response 177518**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 99.10  | 104.13 |
| 106.00 | 93.30  | 98.41  |
| 0.00   | 0.00   | 0.00   |

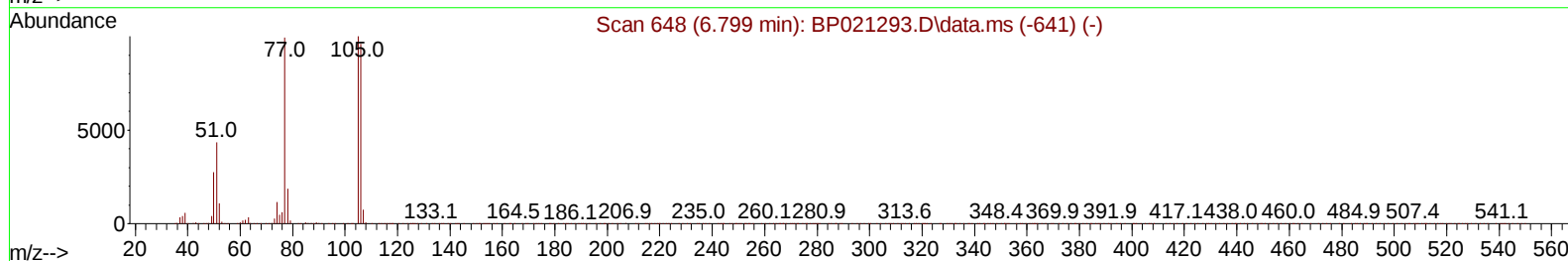
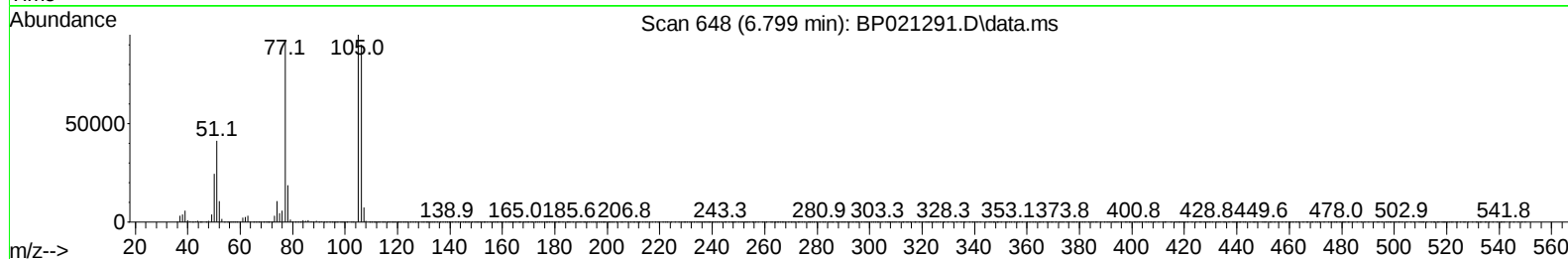
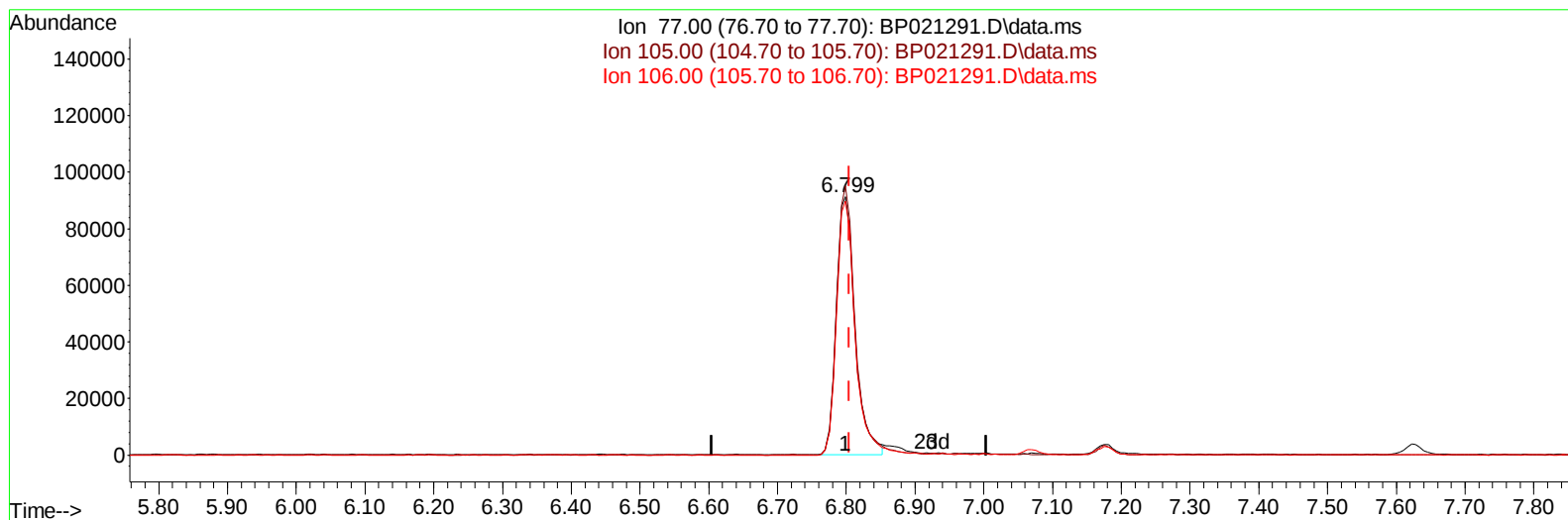
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TIC: BP021291.D\data.ms

**(6) Benzaldehyde**

6.799min (-0.005) 20.37 ng/ul m

response 171274

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 99.10  | 104.13 |
| 106.00 | 93.30  | 98.41  |
| 0.00   | 0.00   | 0.00   |

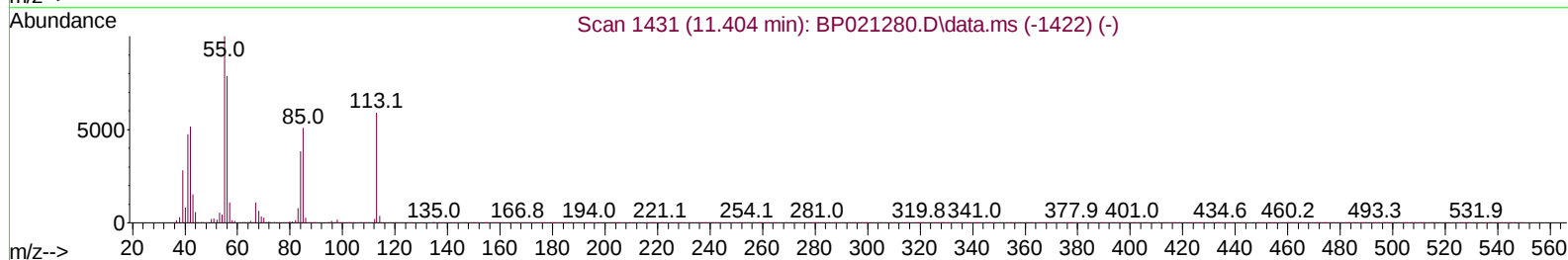
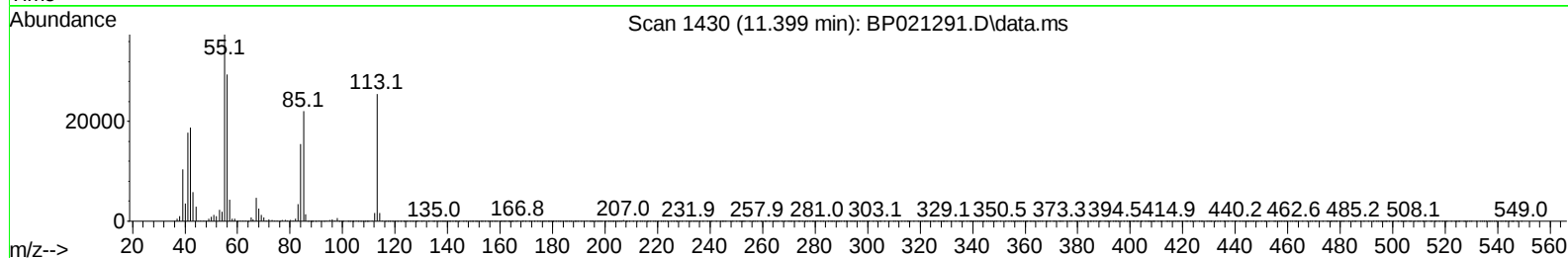
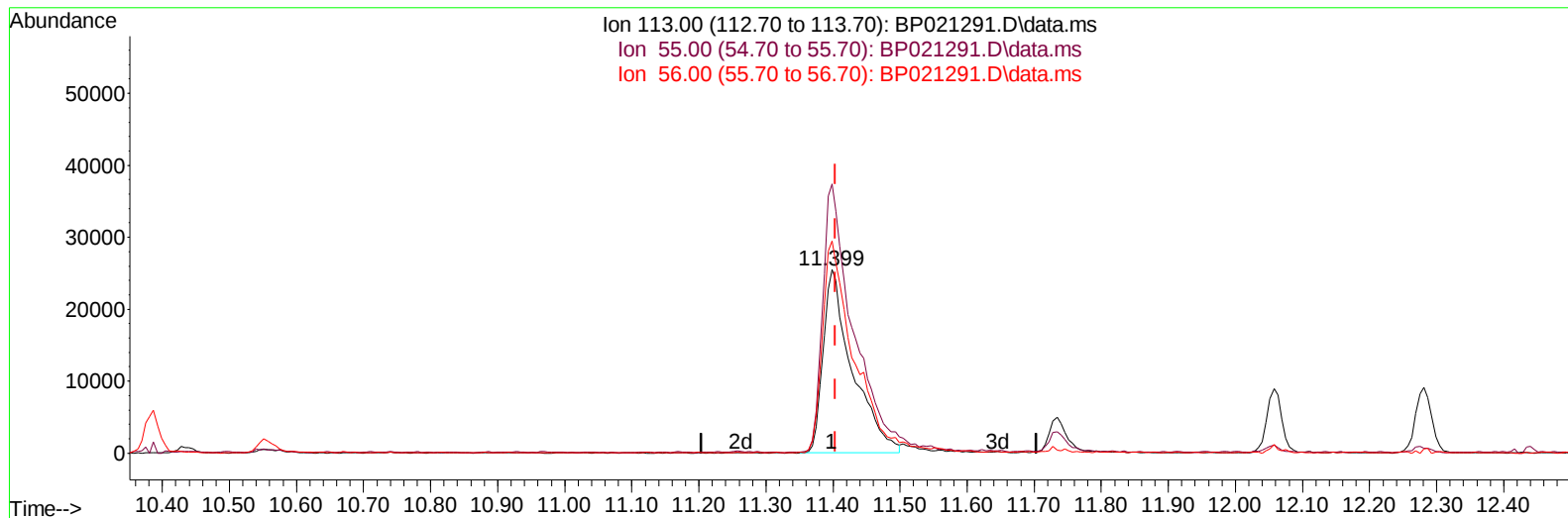
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072924\  
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Misc :  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
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Manual IntegrationsAPPROVED

Quant Time: Aug 01 16:37:43 2024  
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TIC: BP021291.D\data.ms

(34) Caprolactam

11.399min (-0.005) 18.84 ng/ul

response 77376

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 175.30 | 146.72 |
| 56.00  | 131.90 | 115.64 |
| 0.00   | 0.00   | 0.00   |

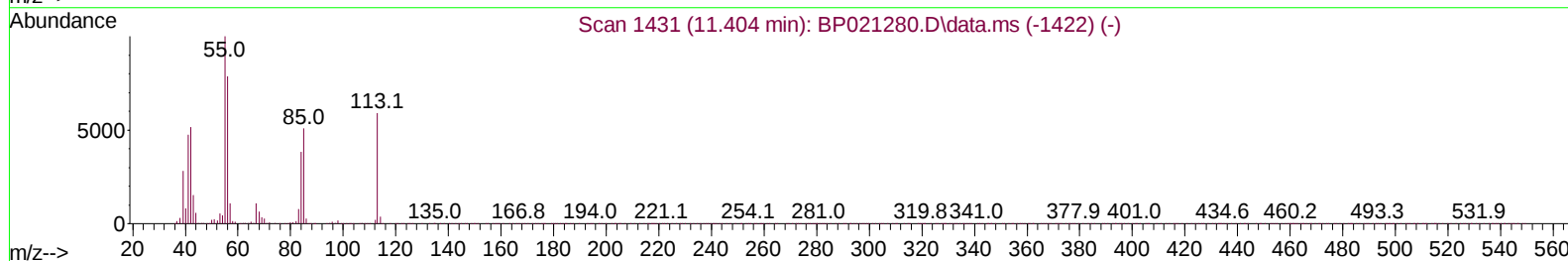
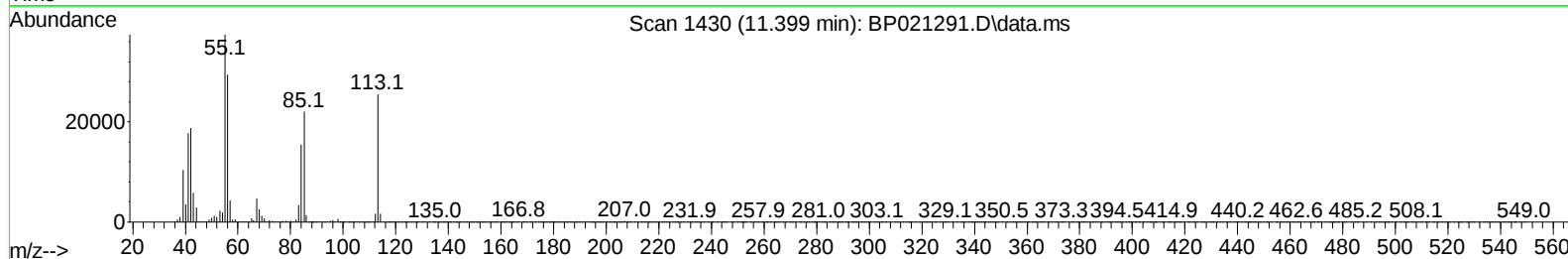
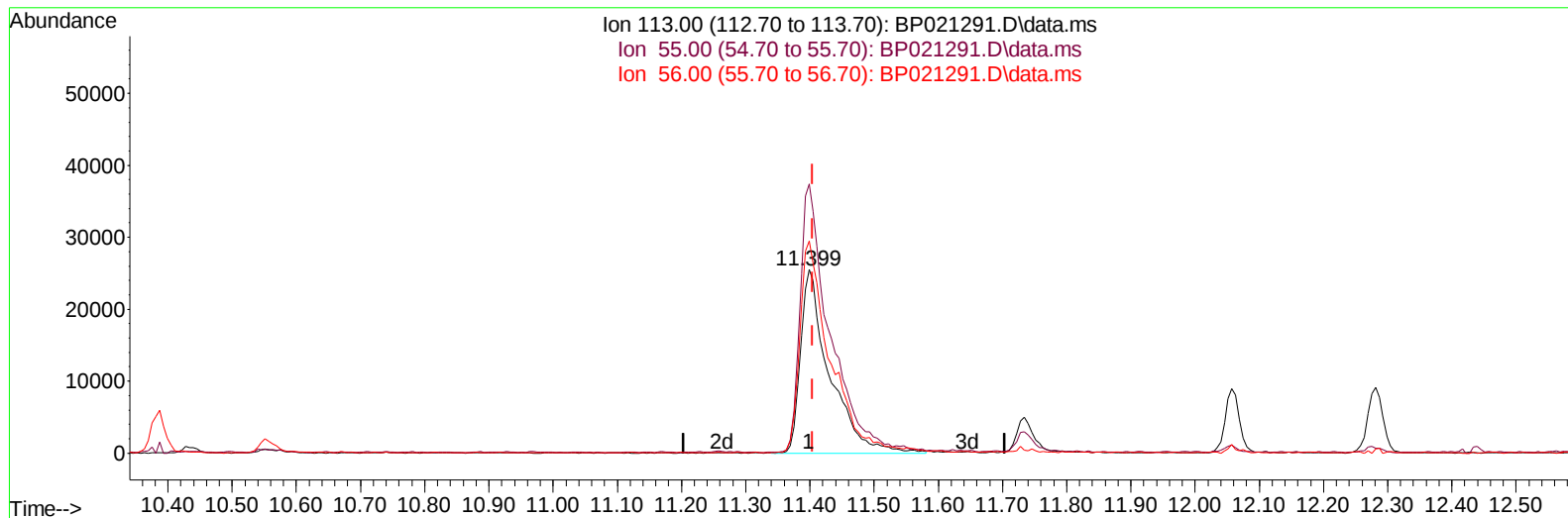
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ALS Vial : 39 Sample Multiplier: 1

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Manual IntegrationsAPPROVED

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TIC: BP021291.D\data.ms

(34) Caprolactam

11.399min (-0.005) 19.57 ng/ul m

response 80383

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 175.30 | 146.72 |
| 56.00  | 131.90 | 115.64 |
| 0.00   | 0.00   | 0.00   |

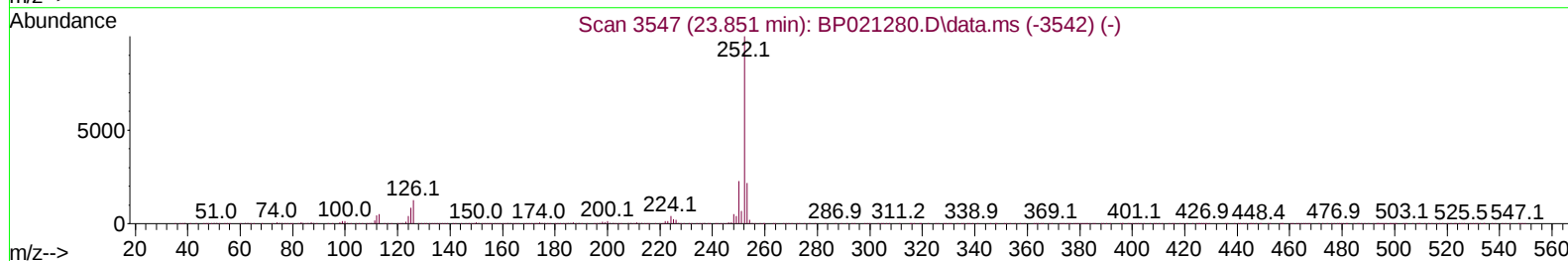
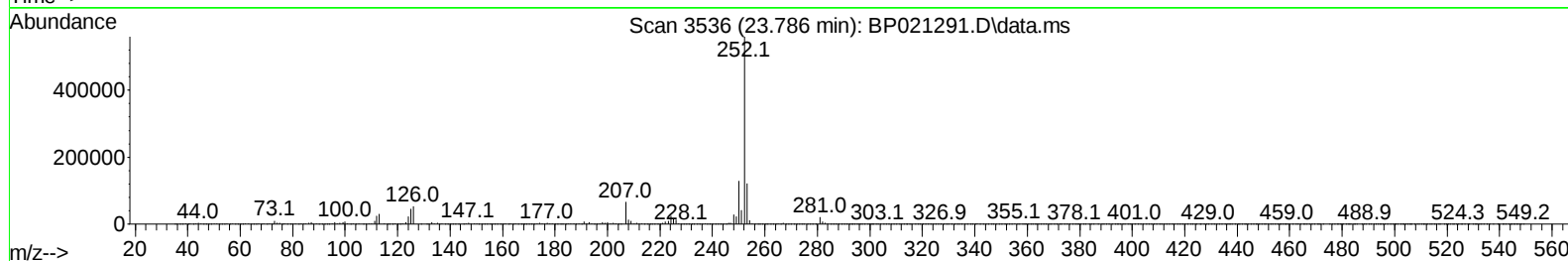
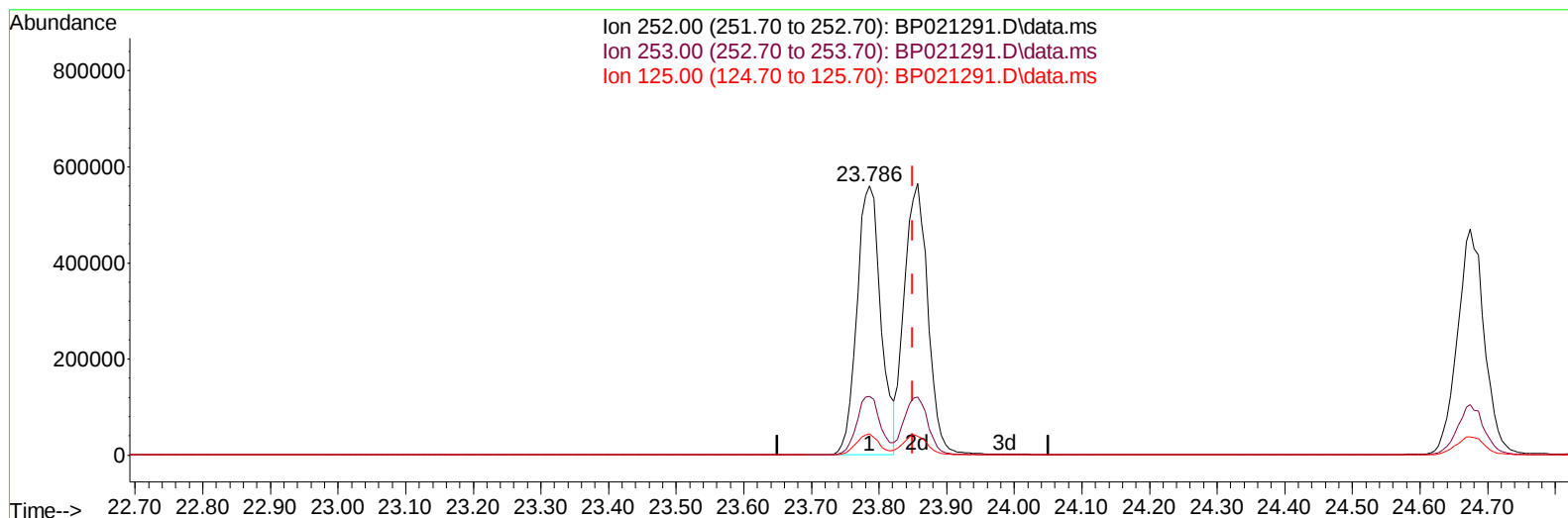
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072924\  
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Sample : SSTDCCC020EC  
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Instrument :  
BNA\_P  
LabSampleId :  
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Manual IntegrationsAPPROVED

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TIC: BP021291.D\data.ms

**(91) Benzo(k)fluoranthene**

**23.786min (-0.064) 18.02 ng/u1**

**response 1383859**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 252.00 | 100.00 | 100.00 |
| 253.00 | 21.80  | 21.81  |
| 125.00 | 9.20   | 8.04   |
| 0.00   | 0.00   | 0.00   |

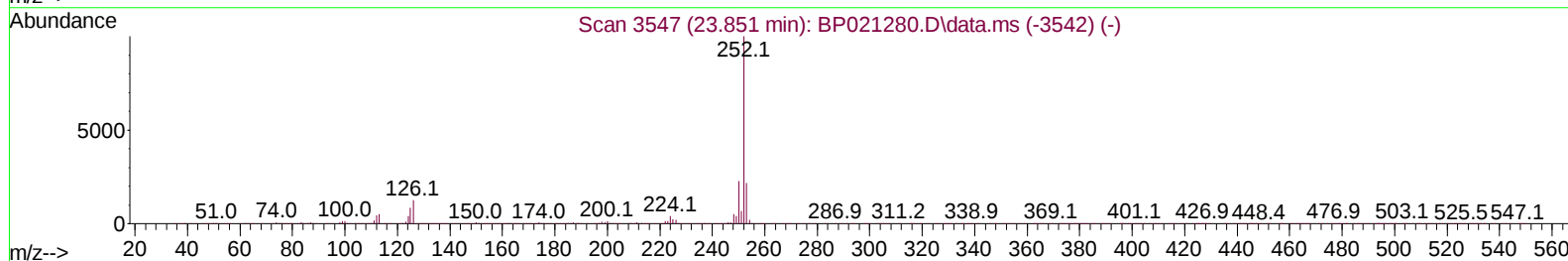
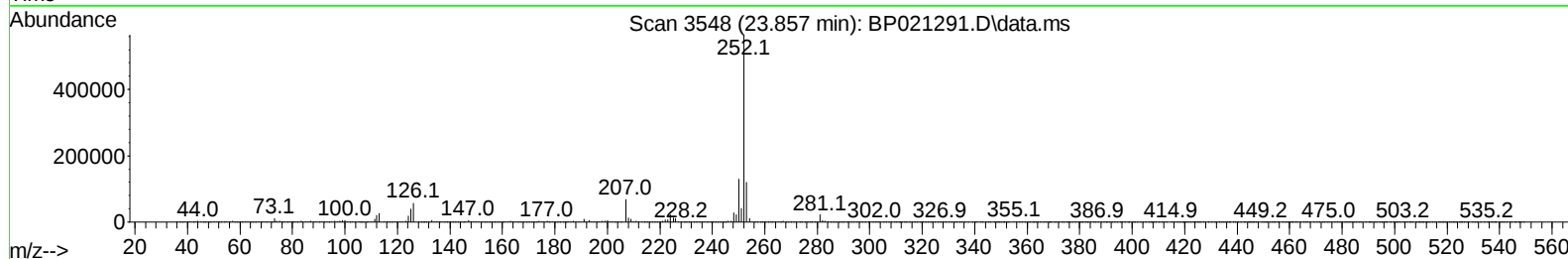
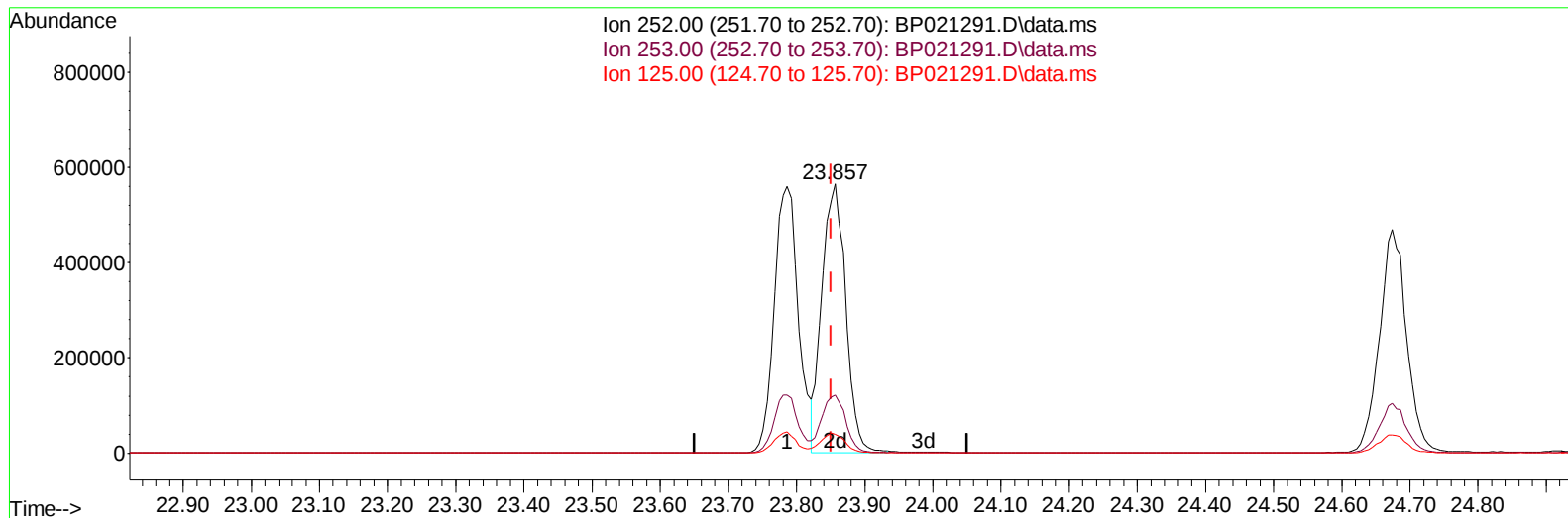
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072924\  
Data File : BP021291.D  
Acq On : 01 Aug 2024 13:24  
Operator : CG/JU  
Sample : SSTDCCC020EC  
Misc :  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020EC

Manual IntegrationsAPPROVED

Quant Time: Aug 01 16:37:43 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 29 12:46:18 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/02/2024  
Supervised By :mohammad ahmed 09/04/2024



TIC: BP021291.D\data.ms

(91) Benzo(k)fluoranthene

23.857min (+ 0.006) 17.83 ng/ul m

response 1369203

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 252.00 | 100.00 | 100.00 |
| 253.00 | 21.80  | 21.48  |
| 125.00 | 9.20   | 7.02#  |
| 0.00   | 0.00   | 0.00   |

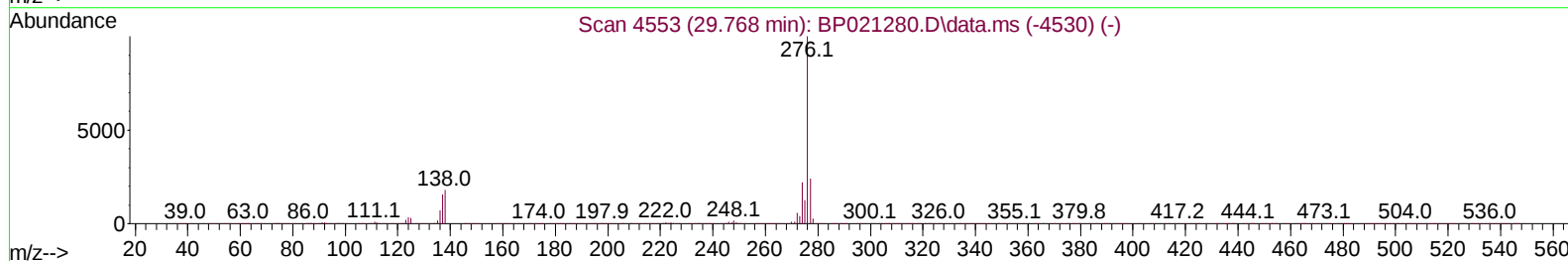
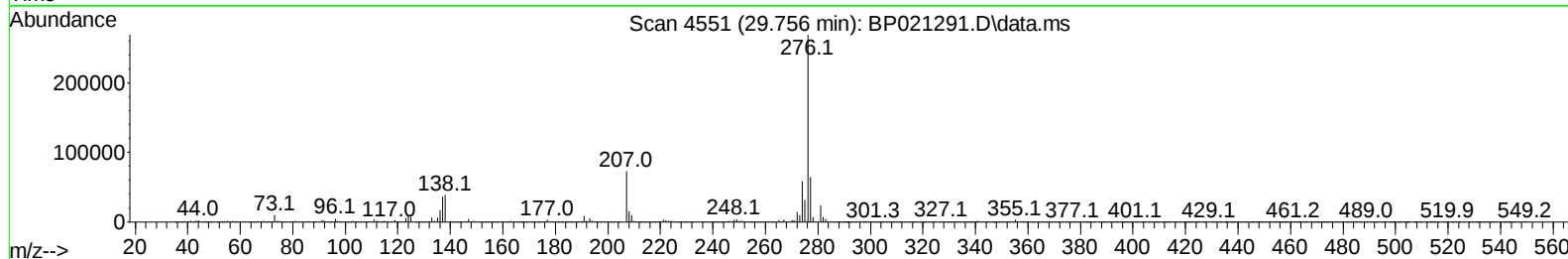
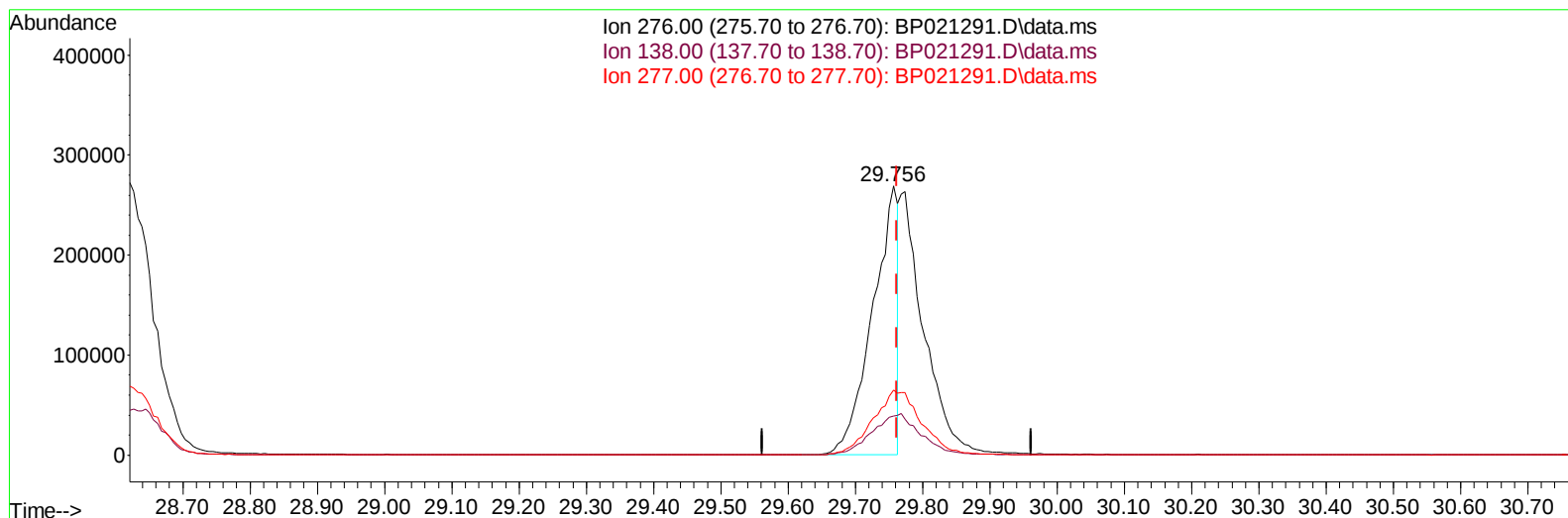
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Instrument :  
 BNA\_P  
 LabSampleId :  
 SSTDCCC020EC

Manual IntegrationsAPPROVED

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TIC: BP021291.D\data.ms

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

29.756min (-0.005) 9.20 ng/u1

response 702045

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 19.10  | 14.47# |
| 277.00 | 22.70  | 24.16  |
| 0.00   | 0.00   | 0.00   |

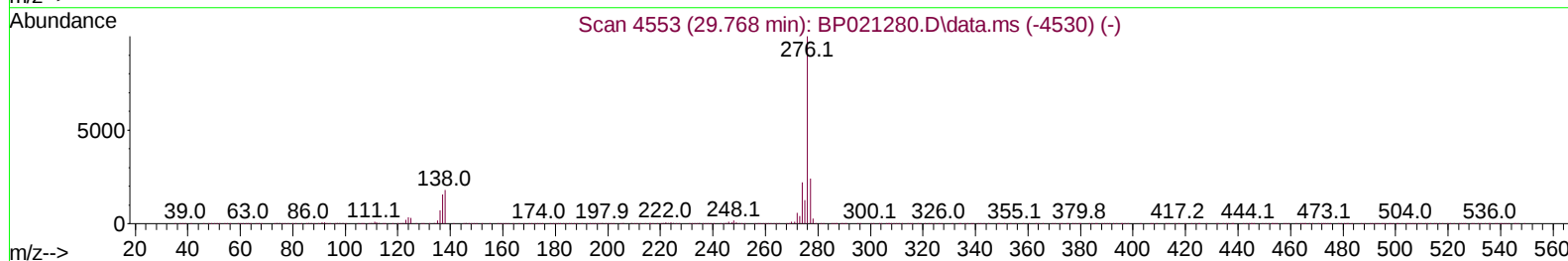
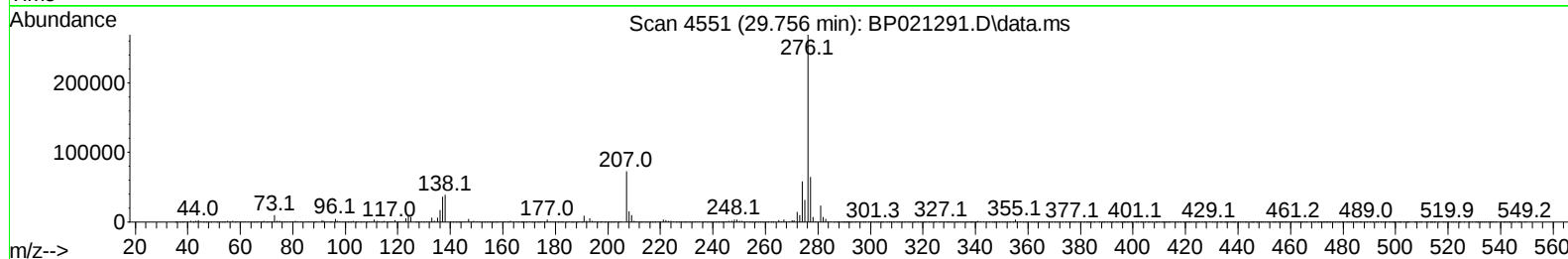
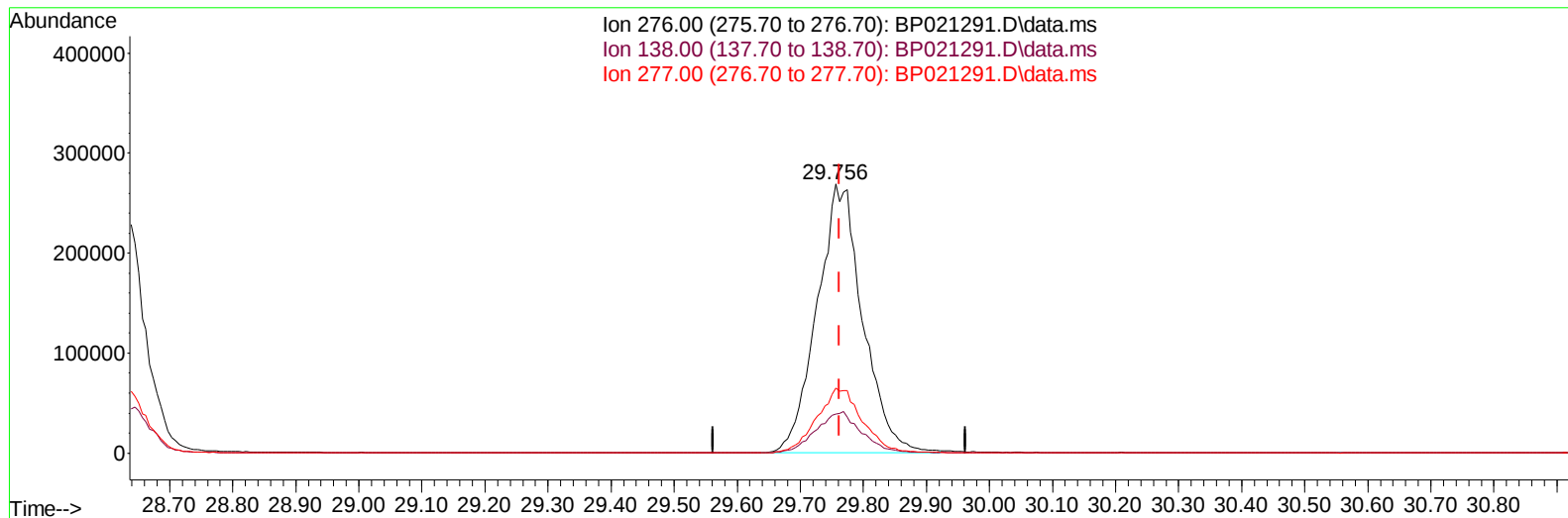
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Data File : BP021291.D  
Acq On : 01 Aug 2024 13:24  
Operator : CG/JU  
Sample : SSTDCCC020EC  
Misc :  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
BNA\_P  
LabSampleId :  
SSTDCCC020EC

Manual IntegrationsAPPROVED

Quant Time: Aug 01 16:37:43 2024  
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Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jul 29 12:46:18 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/02/2024  
Supervised By :mohammad ahmed 09/04/2024



TIC: BP021291.D\data.ms

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

29.756min (-0.005) 17.73 ng/ul m

response 1353276

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 19.10  | 14.47# |
| 277.00 | 22.70  | 24.16  |
| 0.00   | 0.00   | 0.00   |

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 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
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 LabSampleId :  
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Manual IntegrationsAPPROVED

Quant Time: Aug 02 01:53:55 2024  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.628  | 152  | 204593   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.387 | 136  | 841738   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.269 | 164  | 597514   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.092 | 188  | 1344582  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.533 | 240  | 1023572  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.833 | 264  | 1268192  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.164  | 96   | 28729    | 6.390  | ng/uL  | -0.01    |
| 4) Pyridine-d5                | 3.581  | 84   | 205403   | 15.781 | ng/ul  | -0.01    |
| 7) Phenol-d5                  | 6.840  | 99   | 286131   | 17.154 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.981  | 67   | 158624   | 15.692 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.175  | 132  | 252444   | 18.369 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.358  | 113  | 240418   | 17.354 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.793  | 128  | 123208   | 18.845 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.499  | 143  | 145804   | 19.485 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.040 | 165  | 264710   | 19.563 | ng/ul  | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.557 | 131  | 344625   | 19.225 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.687 | 166  | 801029   | 18.442 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.963 | 160  | 920154   | 18.866 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.581 | 143  | 149591   | 24.205 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.287 | 176  | 709553   | 19.912 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.457 | 200  | 142080   | 17.464 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.198 | 188  | 1150735  | 19.269 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.586 | 212  | 1210736  | 19.133 | ng/ul  | 0.01     |
| 92) Benzo(a)pyrene-d12        | 24.598 | 264  | 1153414  | 18.441 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.199  | 88   | 29738    | 6.014  | ng/uL  | 97       |
| 5) Pyridine                   | 3.599  | 79   | 209309   | 15.591 | ng/ul  | 98       |
| 6) Benzaldehyde               | 6.799  | 77   | 171274m  | 20.368 | ng/ul  |          |
| 8) Phenol                     | 6.869  | 94   | 287940   | 17.060 | ng/ul  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.069  | 93   | 228002   | 16.352 | ng/ul  | 95       |
| 12) 2-Chlorophenol            | 7.211  | 128  | 247035   | 17.935 | ng/ul  | 95       |
| 13) 2-Methylphenol            | 8.093  | 108  | 228082   | 17.473 | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.140  | 45   | 283389   | 14.068 | ng/ul# | 92       |
| 16) Acetophenone              | 8.452  | 105  | 370016   | 17.305 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.428  | 70   | 173900   | 15.522 | ng/ul  | 95       |
| 18) 4-Methylphenol            | 8.428  | 108  | 243211   | 17.392 | ng/ul  | 95       |
| 19) Hexachloroethane          | 8.669  | 117  | 106422   | 17.194 | ng/ul  | 92       |
| 22) Nitrobenzene              | 8.828  | 77   | 271931   | 17.363 | ng/ul  | 98       |
| 23) Isophorone                | 9.340  | 82   | 511567   | 16.431 | ng/ul  | 96       |
| 25) 2-Nitrophenol             | 9.534  | 139  | 147817   | 18.940 | ng/ul  | 94       |
| 26) 2,4-Dimethylphenol        | 9.599  | 107  | 277210   | 18.013 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 9.816  | 93   | 311211   | 16.444 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.069 | 162  | 256672   | 19.248 | ng/ul  | 99       |
| 30) Naphthalene               | 10.434 | 128  | 823129   | 18.591 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.575 | 127  | 324928   | 18.983 | ng/ul  | 100      |
| 33) Hexachlorobutadiene       | 10.699 | 225  | 194576   | 19.460 | ng/ul  | 98       |
| 34) Caprolactam               | 11.399 | 113  | 80383m   | 19.568 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.734 | 107  | 283866   | 19.583 | ng/ul  | 96       |

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Manual IntegrationsAPPROVED

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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.057 | 142  | 572096   | 18.823 | ng/ul  | 99       |
| 37) 1-Methylnaphthalene       | 12.281 | 142  | 592174   | 18.945 | ng/ul  | 98       |
| 39) 1,2,4,5-Tetrachloroben... | 12.428 | 216  | 368397   | 18.985 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.387 | 237  | 169245   | 16.673 | ng/ul  | 98       |
| 41) 2,4,6-Trichlorophenol     | 12.698 | 196  | 229265   | 18.698 | ng/ul  | 99       |
| 42) 2,4,5-Trichlorophenol     | 12.798 | 196  | 240877   | 18.567 | ng/ul  | 99       |
| 43) 1,1'-Biphenyl             | 13.081 | 154  | 774364   | 17.733 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.134 | 162  | 638579   | 18.318 | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.369 | 65   | 170426   | 17.678 | ng/ul  | 94       |
| 47) Dimethylphthalate         | 13.734 | 163  | 805696   | 18.281 | ng/ul  | 100      |
| 48) 2,6-Dinitrotoluene        | 13.875 | 165  | 172976   | 19.631 | ng/ul  | 93       |
| 50) Acenaphthylene            | 13.993 | 152  | 1027588  | 18.608 | ng/ul  | 100      |
| 51) 3-Nitroaniline            | 14.222 | 138  | 167371   | 21.828 | ng/ul  | 90       |
| 52) Acenaphthene              | 14.340 | 153  | 688836   | 18.794 | ng/ul  | 97       |
| 53) 2,4-Dinitrophenol         | 14.457 | 184  | 109066   | 23.570 | ng/ul  | 94       |
| 55) 4-Nitrophenol             | 14.592 | 109  | 101973   | 20.285 | ng/ul# | 74       |
| 56) Dibenzofuran              | 14.687 | 168  | 956686   | 19.061 | ng/ul  | 96       |
| 57) 2,4-Dinitrotoluene        | 14.704 | 165  | 258130   | 21.340 | ng/ul# | 81       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.934 | 232  | 225803   | 20.326 | ng/ul  | 94       |
| 59) Diethylphthalate          | 15.116 | 149  | 813084   | 18.460 | ng/ul  | 98       |
| 61) Fluorene                  | 15.345 | 166  | 796977   | 19.459 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.340 | 204  | 436654   | 19.647 | ng/ul  | 94       |
| 63) 4-Nitroaniline            | 15.422 | 138  | 163276   | 25.713 | ng/ul  | 99       |
| 66) 4,6-Dinitro-2-methylph... | 15.475 | 198  | 145382   | 16.851 | ng/ul  | 94       |
| 67) N-Nitrosodiphenylamine    | 15.563 | 169  | 680744   | 17.386 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.257 | 248  | 301114   | 19.106 | ng/ul  | 95       |
| 69) Hexachlorobenzene         | 16.369 | 284  | 365400   | 20.221 | ng/ul  | 93       |
| 70) Atrazine                  | 16.557 | 200  | 277471   | 19.445 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 16.751 | 266  | 230711   | 23.507 | ng/ul  | 96       |
| 72) Phenanthrene              | 17.134 | 178  | 1307408  | 18.744 | ng/ul  | 100      |
| 74) Anthracene                | 17.234 | 178  | 1330516  | 18.711 | ng/ul  | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.051 | 216  | 365177   | 16.969 | ng/uL  | 98       |
| 76) Pentachlorobenzene        | 14.592 | 250  | 394031   | 18.426 | ng/uL  | 99       |
| 77) Carbazole                 | 17.522 | 167  | 1126209  | 19.374 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.098 | 149  | 1445250  | 18.021 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.239 | 202  | 1474625  | 19.296 | ng/ul# | 95       |
| 82) Pyrene                    | 19.616 | 202  | 1454566  | 18.704 | ng/ul  | 96       |
| 83) Butylbenzylphthalate      | 20.551 | 149  | 582284   | 17.971 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.439 | 252  | 422856   | 18.079 | ng/ul  | 98       |
| 85) Benzo(a)anthracene        | 21.516 | 228  | 1293032  | 18.033 | ng/ul  | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.422 | 149  | 818095   | 17.545 | ng/ul# | 97       |
| 87) Chrysene                  | 21.580 | 228  | 1195113  | 17.938 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.674 | 149  | 1226781  | 14.643 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.786 | 252  | 1383859  | 17.982 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.857 | 252  | 1369203m | 17.827 | ng/ul  |          |
| 93) Benzo(a)pyrene            | 24.674 | 252  | 1301792  | 17.900 | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.598 | 276  | 1699821  | 18.138 | ng/ul  | 97       |
| 95) Dibenzo(a,h)anthracene    | 28.651 | 278  | 1421157  | 18.313 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 29.756 | 276  | 1353276m | 17.731 | ng/ul  |          |

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

**Instrument :**  
BNA\_P  
**LabSampleId :**  
SSTDCCC020EC

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 08/02/2024  
Supervised By :mohammad ahmed 09/04/2024

