

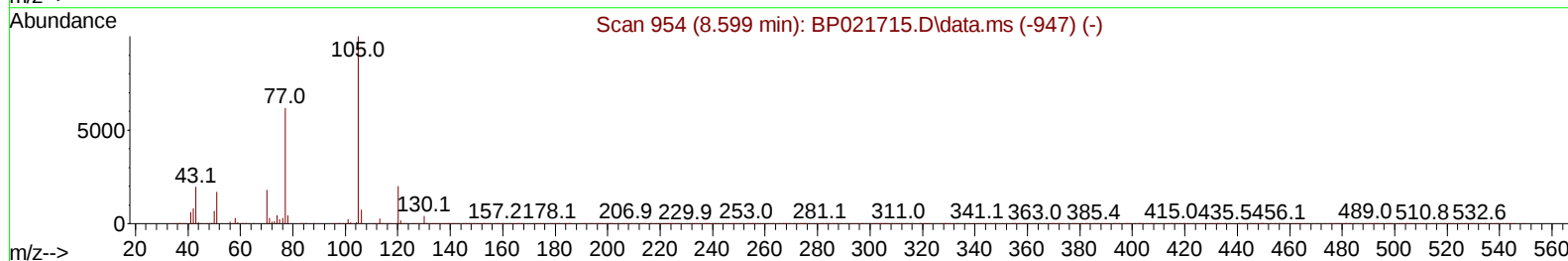
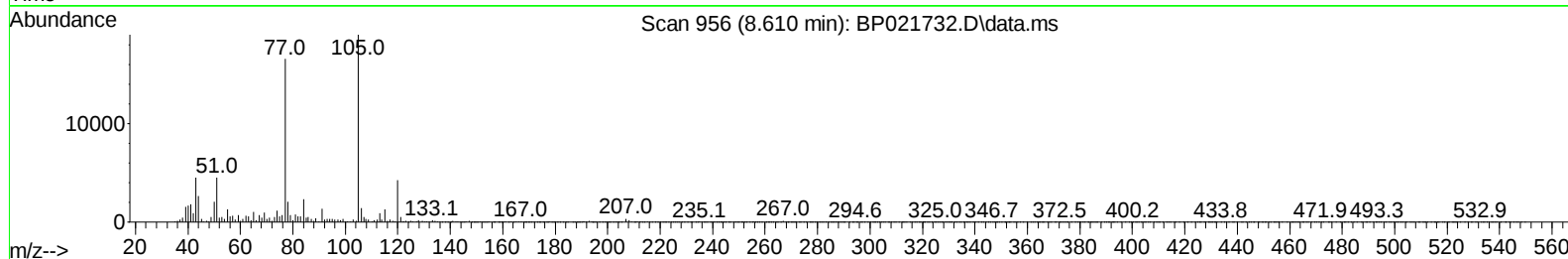
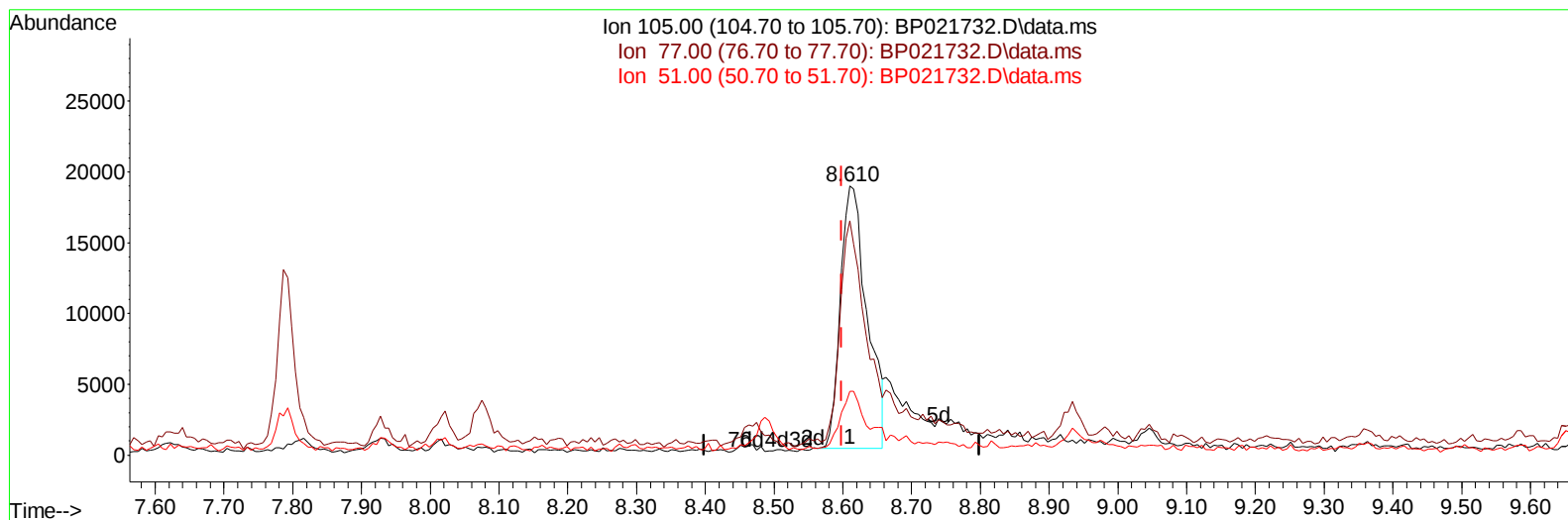
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082924\  
Data File : BP021732.D  
Acq On : 29 Aug 2024 22:34  
Operator : RC/JU  
Sample : P3721-02  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCXY6

Manual IntegrationsAPPROVED

Quant Time: Aug 29 23:59:45 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082324.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 29 23:56:59 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/30/2024  
Supervised By :mohammad ahmed 09/02/2024



TIC: BP021732.D\data.ms

(16) Acetophenone

8.610min (+ 0.011) 0.79 ng/u1

response 49865

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 105.00 | 100.00 | 100.00 |
| 77.00  | 74.30  | 87.20  |
| 51.00  | 20.50  | 23.80  |
| 0.00   | 0.00   | 0.00   |

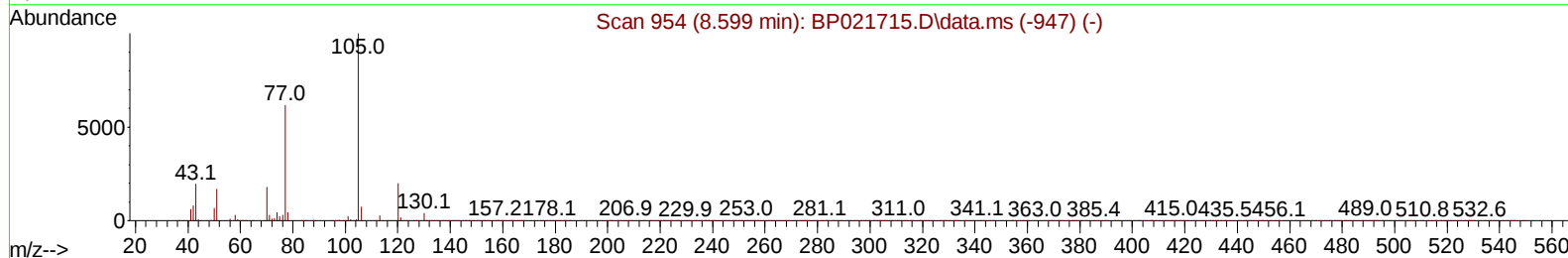
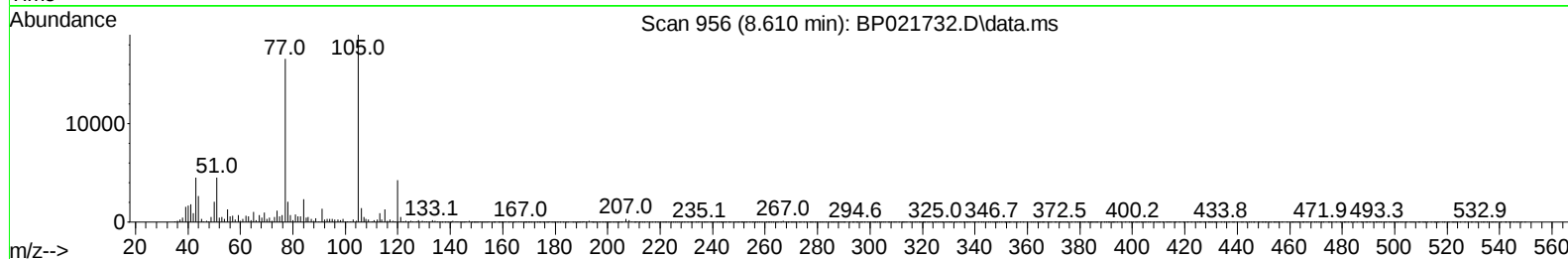
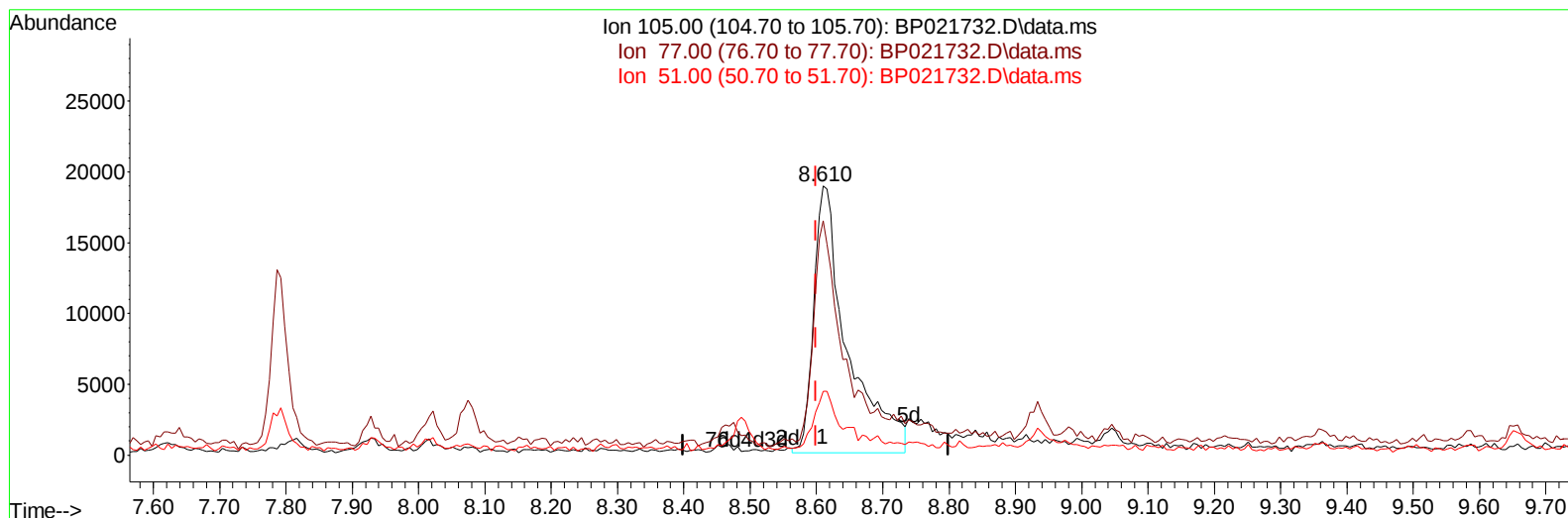
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TIC: BP021732.D\data.ms

(16) Acetophenone

8.610min (+ 0.011) 1.06 ng/ul m

response 66676

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 105.00 | 100.00 | 100.00 |
| 77.00  | 74.30  | 87.20  |
| 51.00  | 20.50  | 23.80  |
| 0.00   | 0.00   | 0.00   |

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 ClientSampleId :  
 DCXY6

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/30/2024  
 Supervised By :mohammad ahmed 09/02/2024

Quant Time: Aug 30 00:08:48 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082324.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 29 23:56:59 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.792  | 152  | 476417   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.575 | 136  | 1978741  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.427 | 164  | 1220648  | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.233 | 188  | 2558631  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.686 | 240  | 2200719  | 20.000 | ng/ul | 0.02     |
| 88) Perylene-d12              | 25.092 | 264  | 2502038  | 20.000 | ng/ul | 0.03     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.275  | 96   | 46154    | 3.216  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 577076   | 13.903 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.957  | 99   | 1203428  | 23.669 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67   | 672296   | 24.297 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.322  | 132  | 800308   | 24.381 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.487  | 113  | 770809   | 19.619 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.934  | 128  | 407900   | 25.622 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.657  | 143  | 431692   | 26.714 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.192 | 165  | 802662   | 25.206 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.704 | 131  | 293889   | 6.344  | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.827 | 166  | 2476580  | 26.516 | ng/ul | -0.01    |
| 49) Acenaphthylene-d8         | 14.116 | 160  | 2816540  | 26.859 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.633 | 143  | 420221   | 23.648 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.439 | 176  | 2101588  | 26.764 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.557 | 200  | 259579   | 18.446 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.327 | 188  | 3203468  | 26.381 | ng/ul | -0.01    |
| 81) Pyrene-d10                | 19.709 | 212  | 3660185  | 29.095 | ng/ul | 0.02     |
| 92) Benzo(a)pyrene-d12        | 24.850 | 264  | 3644907  | 27.320 | ng/ul | 0.01     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 16) Acetophenone              | 8.610  | 105  | 66676m   | 1.059  | ng/ul |          |
| -----                         |        |      |          |        |       |          |

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

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