

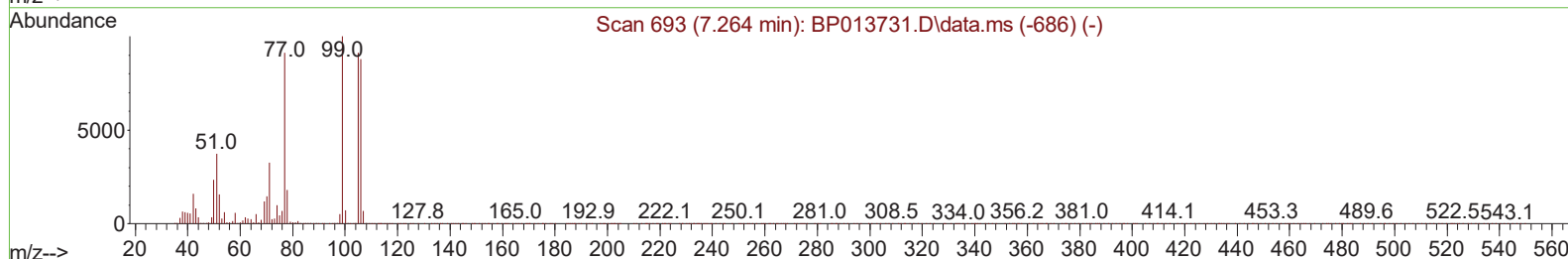
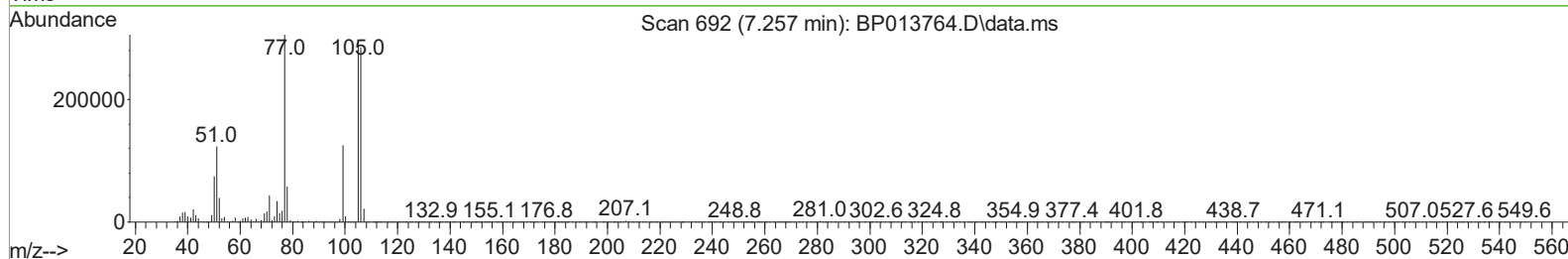
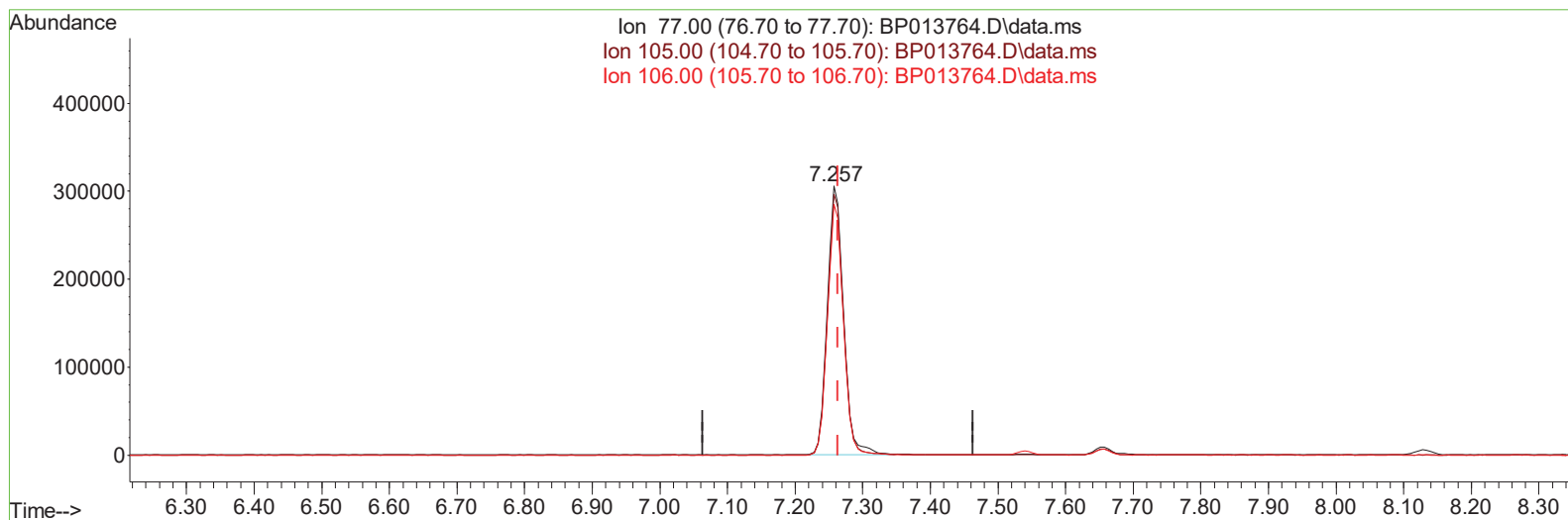
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020223\  
 Data File : BP013763.D  
 Acq On : 04 Feb 2023 05:07  
 Operator : CG/JU  
 Sample : PB150620BS  
 Misc :  
 ALS Vial : 66 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS620

Manual IntegrationsAPPROVED

Quant Time: Feb 04 05:31:49 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP020123.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Feb 03 23:44:43 2023  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/06/2023  
 Supervised By :mohammad ahmed 02/06/2023



TIC: BP013764.D\data.ms

**(6) Benzaldehyde**

7.257min (-0.006) 49.19 ng/ul

response 511865

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 100.80 | 97.11  |
| 106.00 | 93.70  | 93.22  |
| 0.00   | 0.00   | 0.00   |

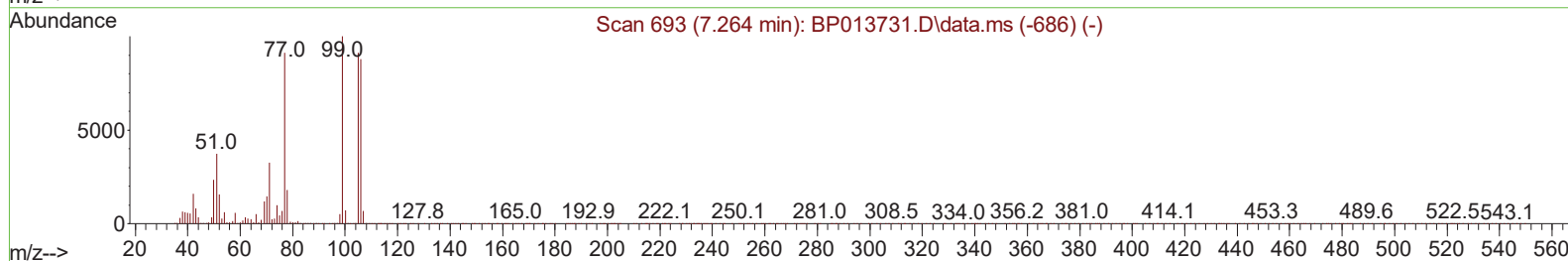
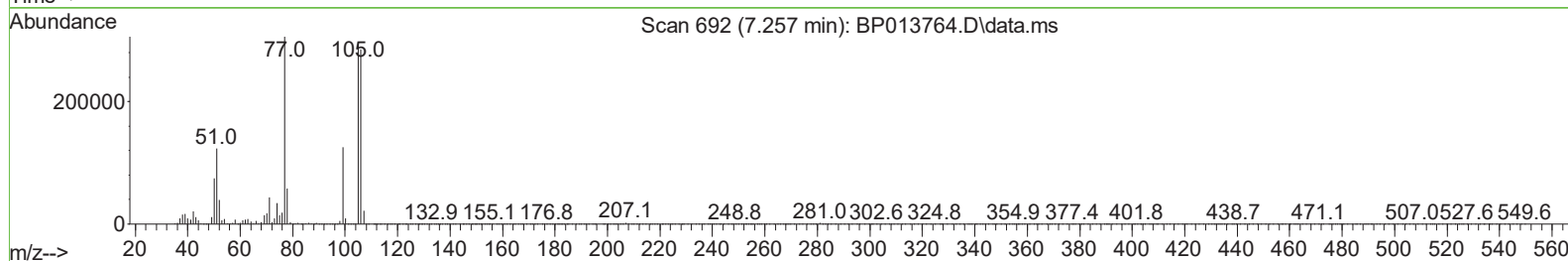
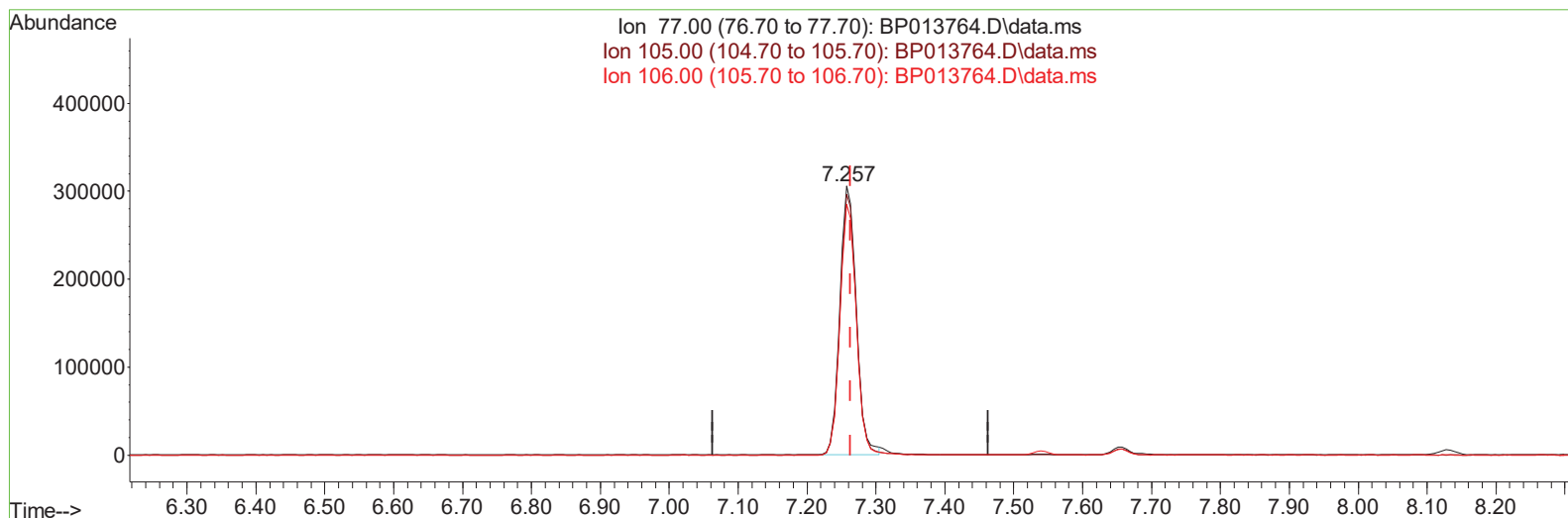
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020223\  
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TIC: BP013764.D\data.ms

**(6) Benzaldehyde**

7.257min (-0.006) 48.60 ng/ul m

response 505720

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 100.80 | 97.11  |
| 106.00 | 93.70  | 93.22  |
| 0.00   | 0.00   | 0.00   |

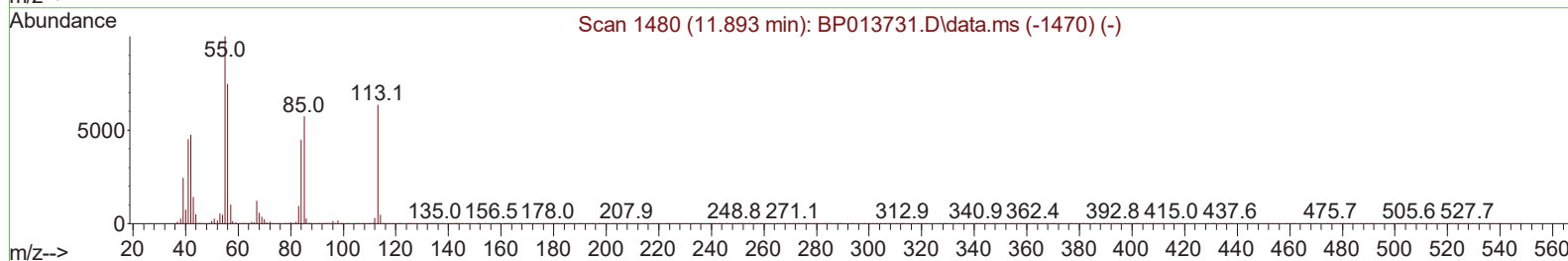
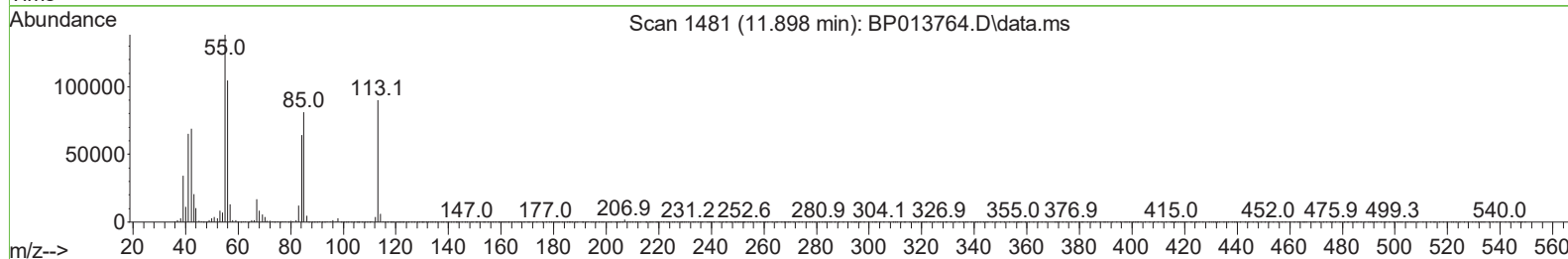
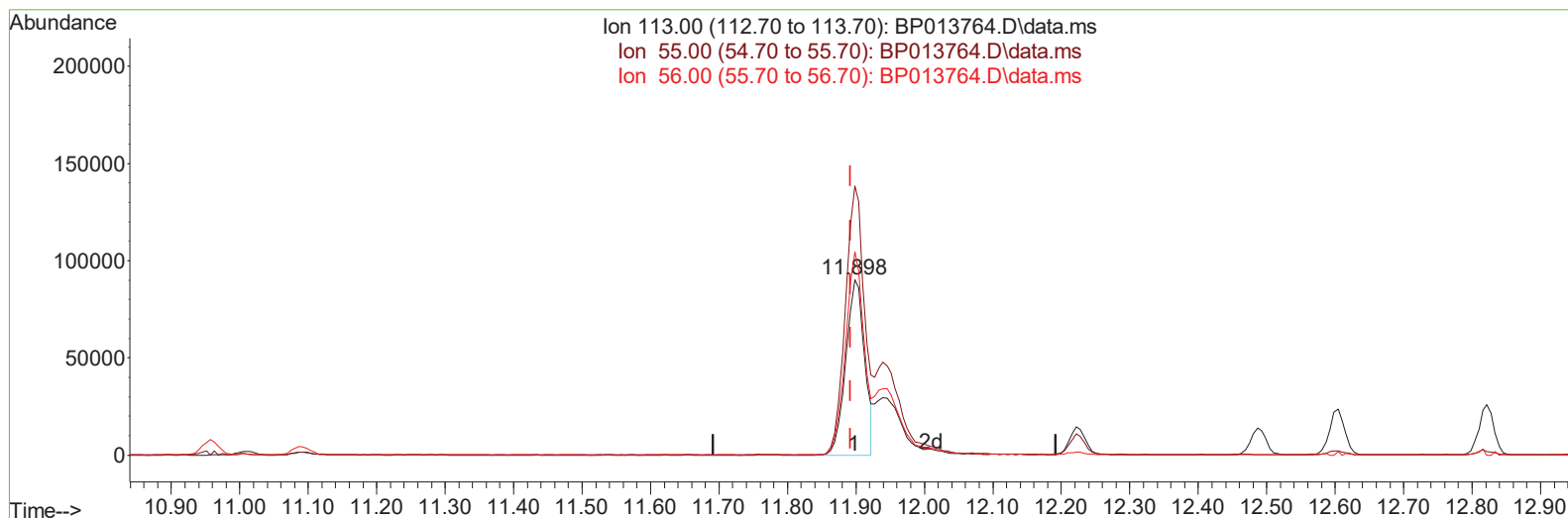
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020223\  
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TIC: BP013764.D\data.ms

**(34) Caprolactam**

11.898min (+ 0.006) 23.34 ng/ul

response 171947

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 160.50 | 153.37 |
| 56.00  | 115.20 | 115.90 |
| 0.00   | 0.00   | 0.00   |

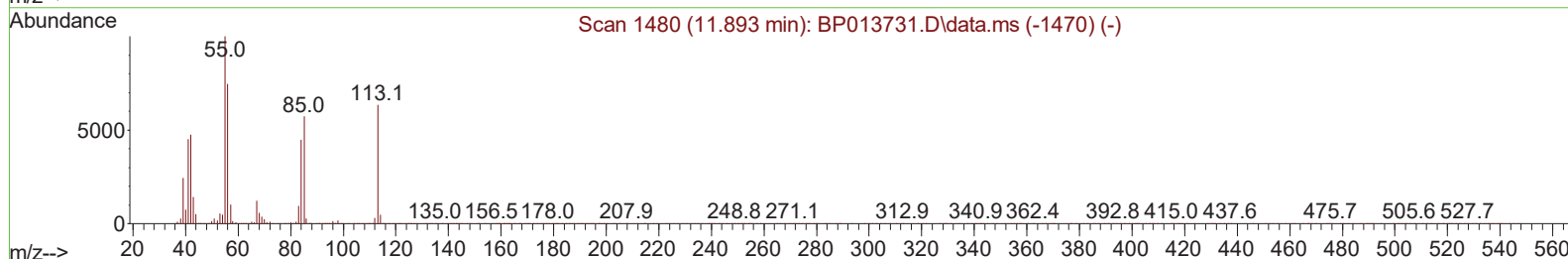
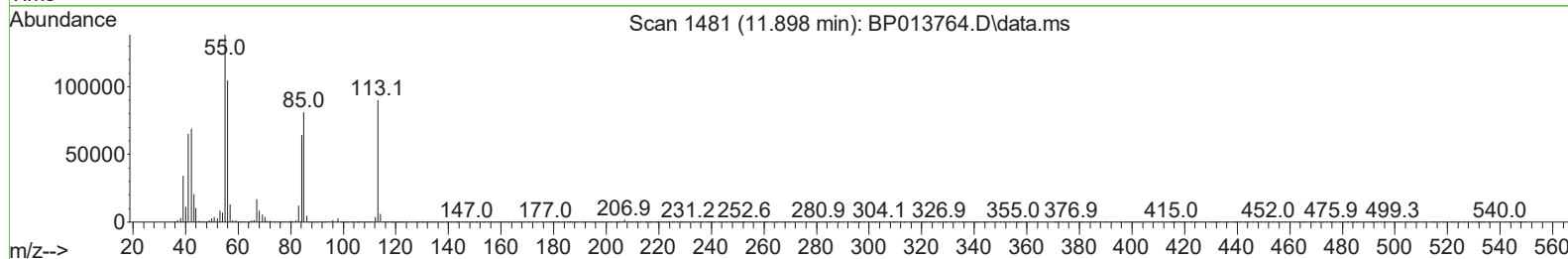
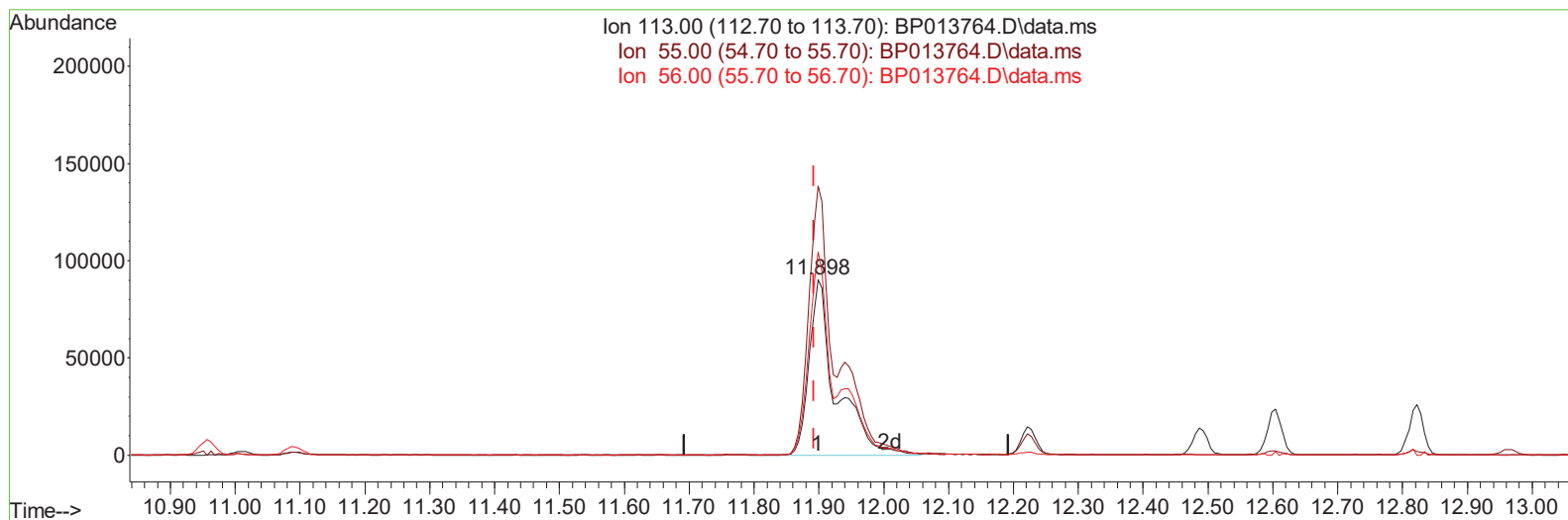
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020223\  
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 Operator : CG/JU  
 Sample : PB150620BS  
 Misc :  
 ALS Vial : 66 Sample Multiplier: 1

Instrument :  
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Manual IntegrationsAPPROVED

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TIC: BP013764.D\data.ms

(34) Caprolactam

11.898min (+ 0.006) 34.93 ng/ul m

response 257330

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 160.50 | 153.37 |
| 56.00  | 115.20 | 115.90 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020223\  
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 ALS Vial : 66 Sample Multiplier: 1

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Manual IntegrationsAPPROVED

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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.128  | 152  | 295296   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.957 | 136  | 1294871  | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.763 | 164  | 932216   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.516 | 188  | 2322774  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.598 | 240  | 2488577  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.162 | 264  | 2363467  | 20.000 | ng/ul  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.475  | 96   | 39467    | 5.667  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.905  | 84   | 592707   | 29.659 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.275  | 99   | 853675   | 32.893 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.446  | 67   | 489960   | 32.189 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.657  | 132  | 625259   | 32.865 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.834  | 113  | 677651   | 32.753 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.304  | 128  | 300134   | 36.334 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.028 | 143  | 355224   | 36.900 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.569 | 165  | 749622   | 33.299 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.092 | 131  | 842267   | 28.329 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.169 | 166  | 2304053  | 31.355 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.463 | 160  | 2535641  | 31.232 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.951 | 143  | 390525   | 31.530 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.757 | 176  | 2114605  | 33.386 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.869 | 200  | 330091   | 25.359 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.616 | 188  | 3441161  | 31.783 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.833 | 212  | 4293601  | 30.934 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.998 | 264  | 4011372  | 33.247 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.510  | 88   | 87613    | 12.071 | ng/uL  | 97       |
| 5) Pyridine                   | 3.928  | 79   | 633566   | 31.485 | ng/ul  | 97       |
| 6) Benzaldehyde               | 7.257  | 77   | 505720m  | 48.598 | ng/ul  |          |
| 8) Phenol                     | 7.304  | 94   | 929917   | 35.012 | ng/ul  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.540  | 93   | 713693   | 33.439 | ng/ul  | 96       |
| 12) 2-Chlorophenol            | 7.687  | 128  | 671923   | 34.861 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.569  | 108  | 693729   | 34.209 | ng/ul  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.651  | 45   | 849565   | 33.620 | ng/ul  | 98       |
| 16) Acetophenone              | 8.963  | 105  | 1097305  | 34.021 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.945  | 70   | 560688   | 34.702 | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.898  | 108  | 771400   | 35.025 | ng/ul  | 99       |
| 19) Hexachloroethane          | 9.222  | 117  | 258389   | 34.249 | ng/ul  | 98       |
| 22) Nitrobenzene              | 9.345  | 77   | 814278   | 36.497 | ng/ul  | 98       |
| 23) Isophorone                | 9.875  | 82   | 1658689  | 32.474 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 10.063 | 139  | 399945   | 37.549 | ng/ul  | 97       |
| 26) 2,4-Dimethylphenol        | 10.116 | 107  | 809734   | 33.647 | ng/ul  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 10.351 | 93   | 1016650  | 32.842 | ng/ul  | 100      |
| 29) 2,4-Dichlorophenol        | 10.598 | 162  | 761255   | 34.942 | ng/ul  | 99       |
| 30) Naphthalene               | 11.010 | 128  | 2234588  | 32.652 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 11.116 | 127  | 910271   | 30.421 | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 11.292 | 225  | 516665   | 33.626 | ng/ul  | 97       |
| 34) Caprolactam               | 11.898 | 113  | 257330m  | 34.933 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.222 | 107  | 801859   | 34.363 | ng/ul  | 99       |

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 ALS Vial : 66 Sample Multiplier: 1

Instrument :  
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 ClientSampleId :  
 SLCS620

## Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 02/06/2023

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| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.604 | 142  | 1608856  | 33.026 | ng/ul | 99       |
| 37) 1-Methylnaphthalene       | 12.822 | 142  | 1643415  | 33.403 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.969 | 216  | 1047745  | 34.025 | ng/ul | 99       |
| 40) Hexachlorocyclopentadiene | 12.945 | 237  | 493712   | 28.516 | ng/ul | 99       |
| 41) 2,4,6-Trichlorophenol     | 13.204 | 196  | 698864   | 34.260 | ng/ul | 100      |
| 42) 2,4,5-Trichlorophenol     | 13.275 | 196  | 766301   | 34.470 | ng/ul | 99       |
| 43) 1,1'-Biphenyl             | 13.598 | 154  | 2278797  | 32.286 | ng/ul | 100      |
| 44) 2-Chloronaphthalene       | 13.645 | 162  | 1827690  | 32.731 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.845 | 65   | 464224   | 43.673 | ng/ul | 98       |
| 47) Dimethylphthalate         | 14.216 | 163  | 2384302  | 32.935 | ng/ul | 99       |
| 48) 2,6-Dinitrotoluene        | 14.339 | 165  | 464769   | 45.048 | ng/ul | 96       |
| 50) Acenaphthylene            | 14.486 | 152  | 2810001  | 32.823 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.669 | 138  | 437856   | 40.152 | ng/ul | 100      |
| 52) Acenaphthene              | 14.827 | 153  | 1936167  | 32.910 | ng/ul | 100      |
| 53) 2,4-Dinitrophenol         | 14.869 | 184  | 139174   | 17.368 | ng/ul | 99       |
| 55) 4-Nitrophenol             | 14.963 | 109  | 307802   | 32.600 | ng/ul | 98       |
| 56) Dibenzofuran              | 15.163 | 168  | 2825526  | 33.064 | ng/ul | 99       |
| 57) 2,4-Dinitrotoluene        | 15.122 | 165  | 689880   | 42.573 | ng/ul | 98       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.386 | 232  | 710192   | 33.991 | ng/ul | 97       |
| 59) Diethylphthalate          | 15.574 | 149  | 2308559  | 33.217 | ng/ul | 99       |
| 61) Fluorene                  | 15.810 | 166  | 2428974  | 34.926 | ng/ul | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.798 | 204  | 1349022  | 35.604 | ng/ul | 98       |
| 63) 4-Nitroaniline            | 15.833 | 138  | 494630   | 48.442 | ng/ul | 98       |
| 66) 4,6-Dinitro-2-methylph... | 15.880 | 198  | 336964   | 26.012 | ng/ul | 99       |
| 67) N-Nitrosodiphenylamine    | 16.016 | 169  | 2081242  | 33.196 | ng/ul | 100      |
| 68) 4-Bromophenyl-phenylether | 16.698 | 248  | 844505   | 31.956 | ng/ul | 98       |
| 69) Hexachlorobenzene         | 16.816 | 284  | 1003289  | 32.360 | ng/ul | 99       |
| 70) Atrazine                  | 16.963 | 200  | 826320   | 32.748 | ng/ul | 98       |
| 71) Pentachlorophenol         | 17.163 | 266  | 588839   | 29.756 | ng/ul | 98       |
| 72) Phenanthrene              | 17.563 | 178  | 4097895  | 33.399 | ng/ul | 100      |
| 74) Anthracene                | 17.651 | 178  | 4150481  | 33.399 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.569 | 216  | 1056155  | 31.916 | ng/uL | 99       |
| 76) Pentachlorobenzene        | 15.080 | 250  | 1083516  | 31.939 | ng/uL | 99       |
| 77) Carbazole                 | 17.916 | 167  | 3715471  | 34.463 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 18.445 | 149  | 4184254  | 32.848 | ng/ul | 100      |
| 80) Fluoranthene              | 19.510 | 202  | 5405887  | 33.415 | ng/ul | 99       |
| 82) Pyrene                    | 19.862 | 202  | 5378187  | 32.516 | ng/ul | 99       |
| 83) Butylbenzylphthalate      | 20.715 | 149  | 1810949  | 38.108 | ng/ul | 97       |
| 84) 3,3'-Dichlorobenzidine    | 21.504 | 252  | 1777255  | 31.443 | ng/ul | 100      |
| 85) Benzo(a)anthracene        | 21.580 | 228  | 5861134  | 34.259 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.474 | 149  | 2851639  | 35.000 | ng/ul | 99       |
| 87) Chrysene                  | 21.639 | 228  | 5470389  | 33.962 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.456 | 149  | 4995464  | 40.124 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.374 | 252  | 5839543  | 37.849 | ng/ul | 100      |
| 91) Benzo(k)fluoranthene      | 23.427 | 252  | 5669626  | 38.449 | ng/ul | 99       |
| 93) Benzo(a)pyrene            | 24.051 | 252  | 4654233  | 34.942 | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.880 | 276  | 5477116  | 33.189 | ng/ul | 98       |
| 95) Dibenzo(a,h)anthracene    | 26.892 | 278  | 4599877  | 33.622 | ng/ul | 100      |
| 96) Benzo(g,h,i)perylene      | 27.709 | 276  | 4354315  | 33.352 | ng/ul | 99       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SLCS620

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