

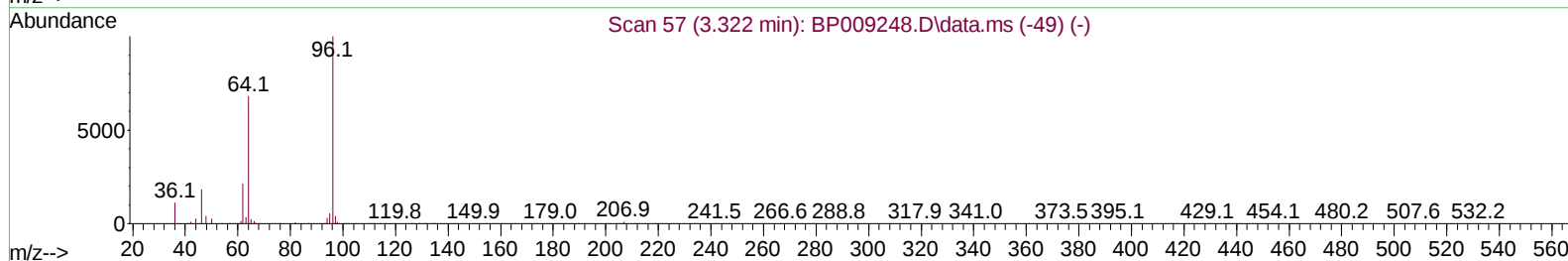
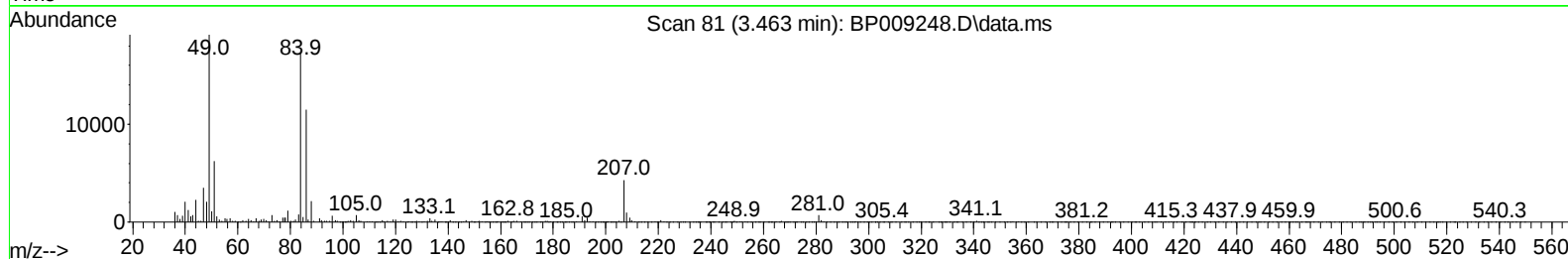
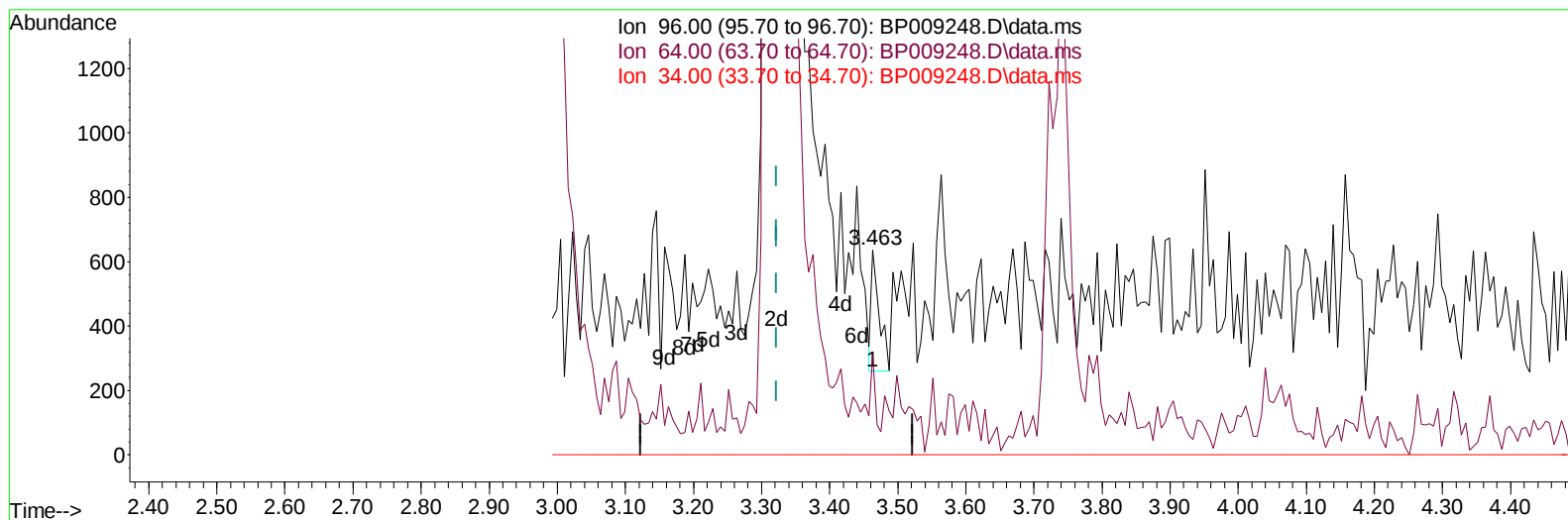
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP030222\  
 Data File : BP009248.D  
 Acq On : 02 Mar 2022 16:39  
 Operator : CG/JU  
 Sample : SST02010  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST020614

Manual IntegrationsAPPROVED

Quant Time: Mar 02 17:37:09 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030222.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Mar 02 17:32:26 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/03/2022  
 Supervised By :mohammad ahmed 03/07/2022



TIC: BP009248.D\data.ms

**(3) 1,4-Dioxane-d8 (S)**

**3.463min (+ 0.141) 0.04 ng/uL**

**response 304**

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 55.40  | 48.98  |
| 34.00 | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

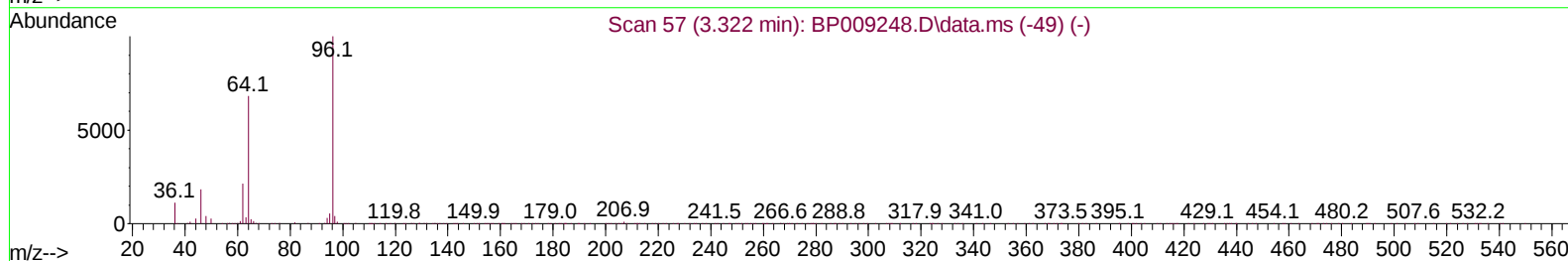
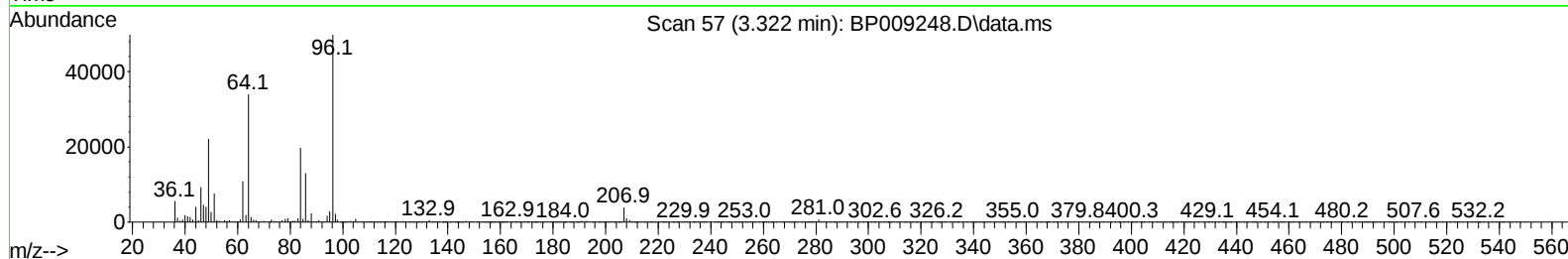
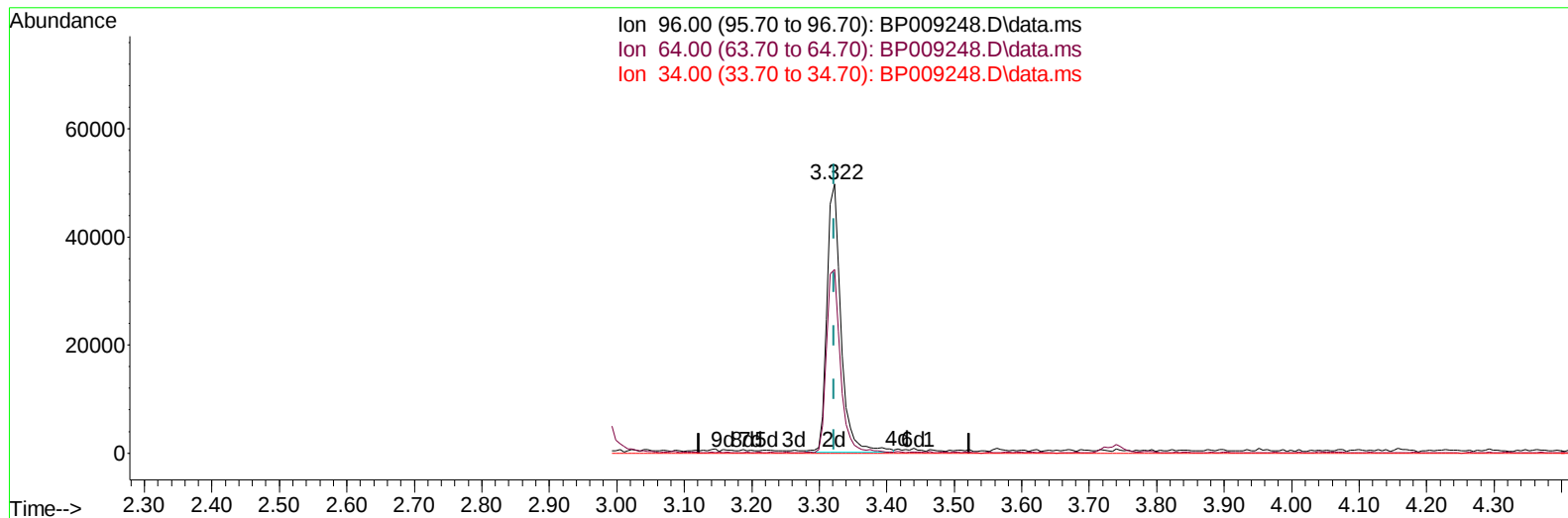
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP030222\  
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Manual IntegrationsAPPROVED

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TIC: BP009248.D\data.ms

**(3) 1,4-Dioxane-d8 (S)**

3.322min ( 0.000) 8.40 ng/uL m

response 71704

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 55.40  | 68.26# |
| 34.00 | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

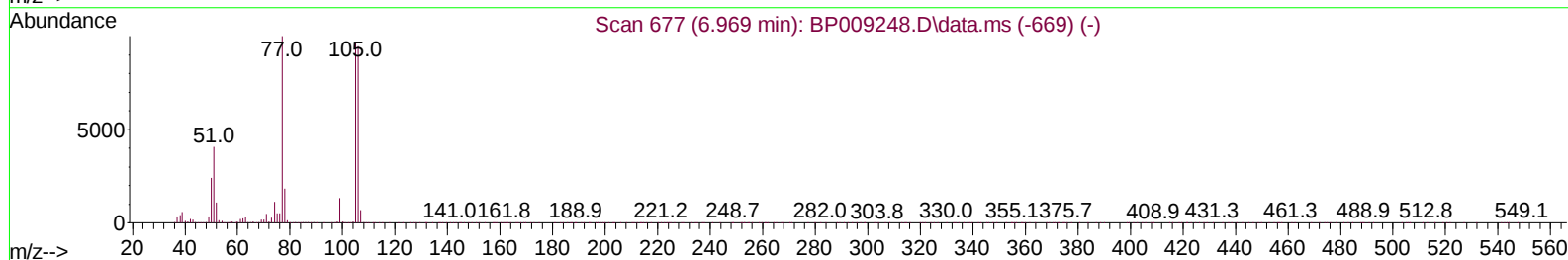
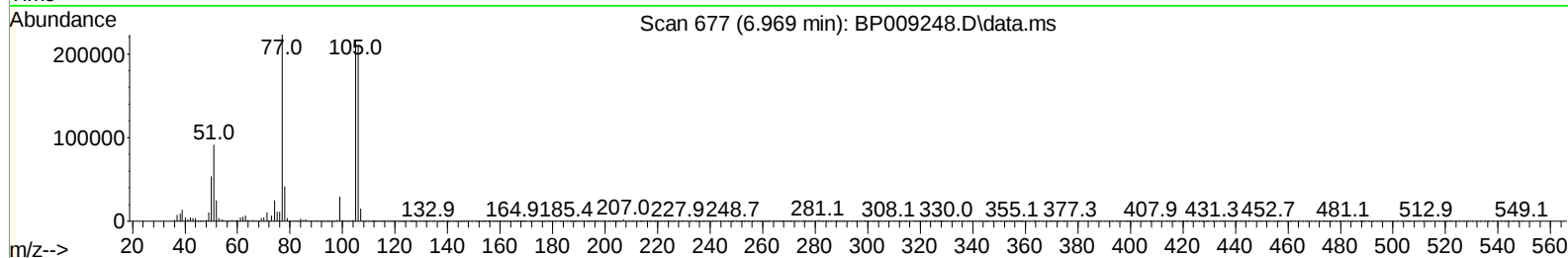
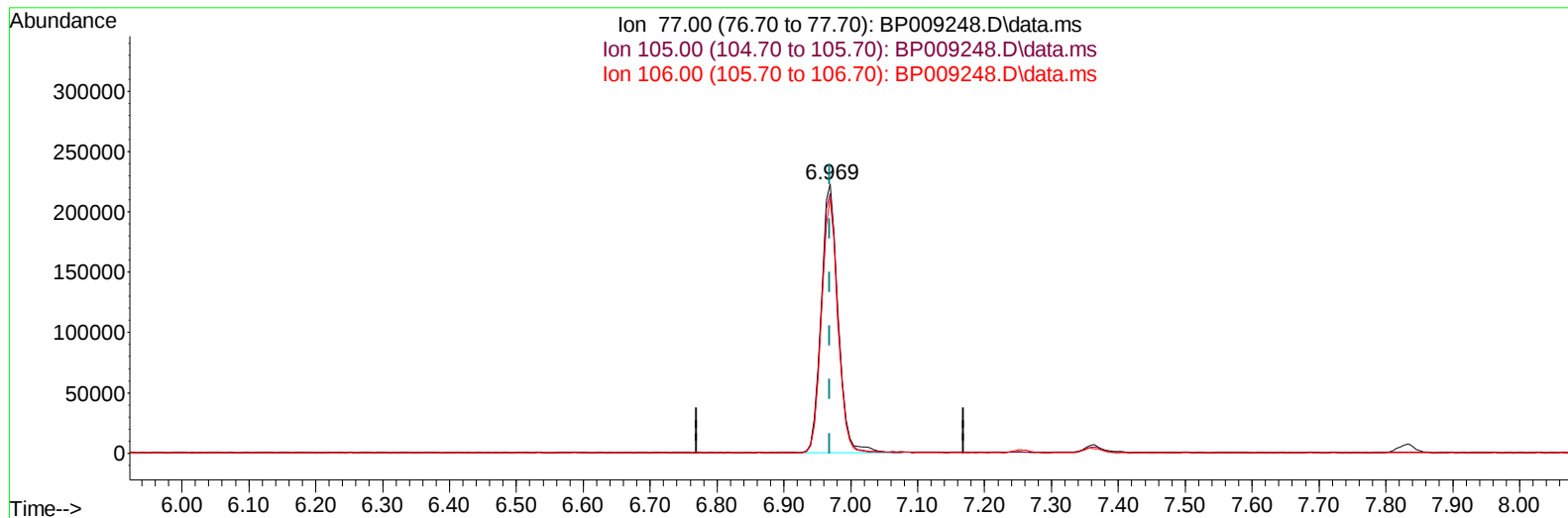
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP030222\  
Data File : BP009248.D  
Acq On : 02 Mar 2022 16:39  
Operator : CG/JU  
Sample : SSTD02010  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD020614

Manual IntegrationsAPPROVED

Quant Time: Mar 02 17:37:09 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030222.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Mar 02 17:32:26 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/03/2022  
Supervised By :mohammad ahmed 03/07/2022



TIC: BP009248.D\data.ms

(6) Benzaldehyde

6.969min ( 0.000) 23.31 ng/u1

response 390600

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 96.20  | 96.50  |
| 106.00 | 96.20  | 94.66  |
| 0.00   | 0.00   | 0.00   |

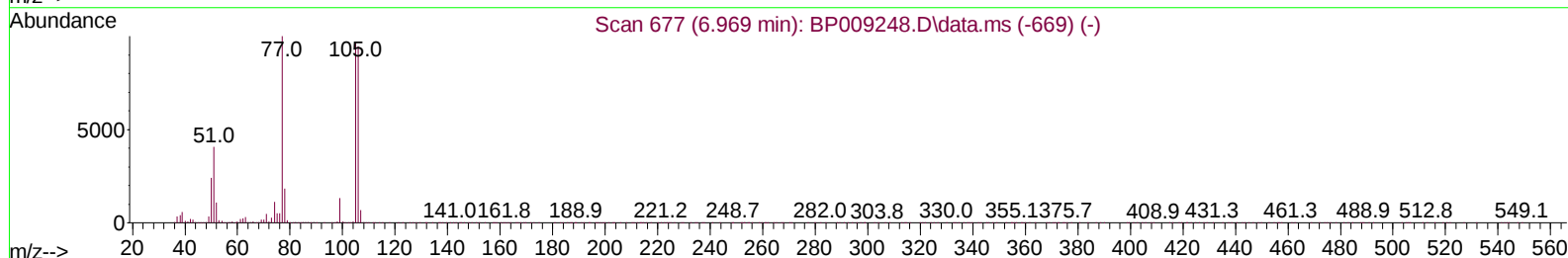
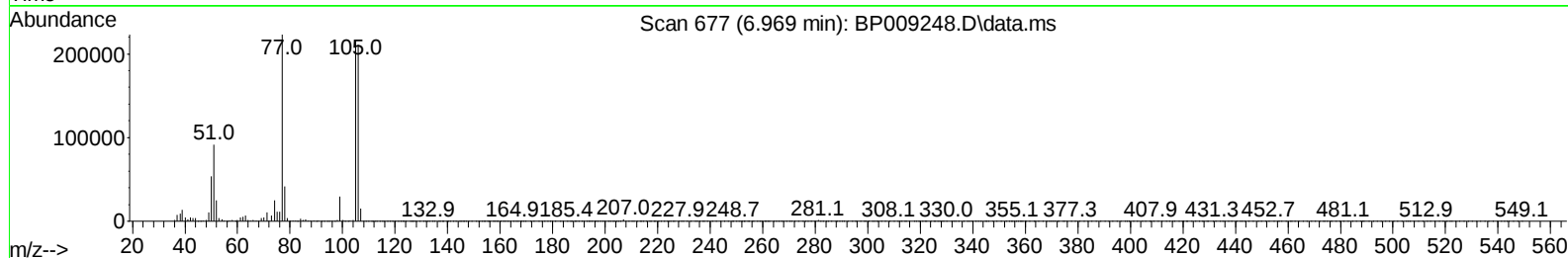
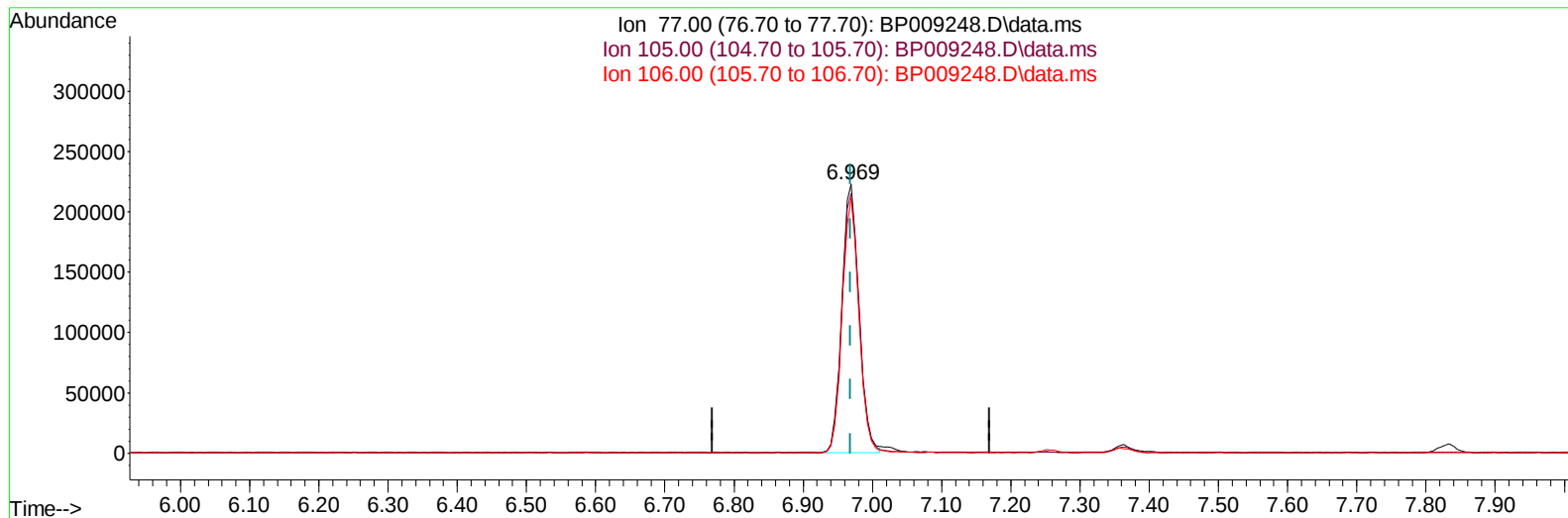
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP030222\  
 Data File : BP009248.D  
 Acq On : 02 Mar 2022 16:39  
 Operator : CG/JU  
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 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
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Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 03/07/2022



TIC: BP009248.D\data.ms

**(6) Benzaldehyde**

6.969min ( 0.000) 22.91 ng/ul m

response 383958

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 77.00  | 100.00 | 100.00 |
| 105.00 | 96.20  | 96.50  |
| 106.00 | 96.20  | 94.66  |
| 0.00   | 0.00   | 0.00   |

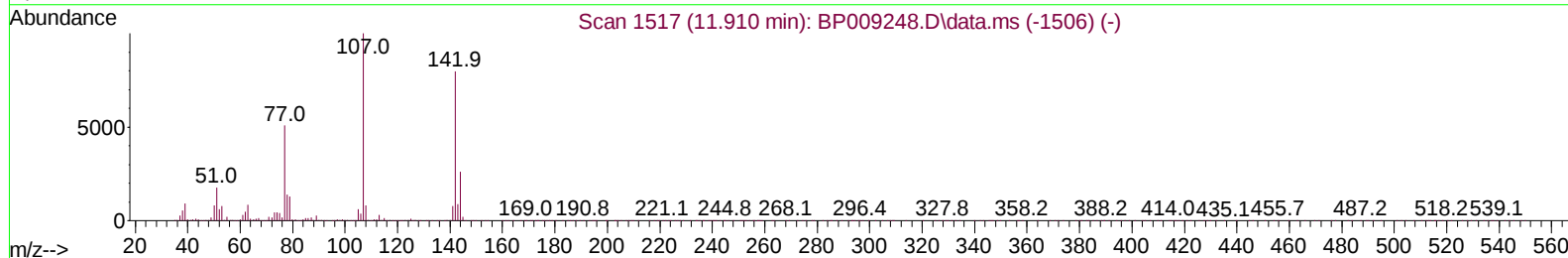
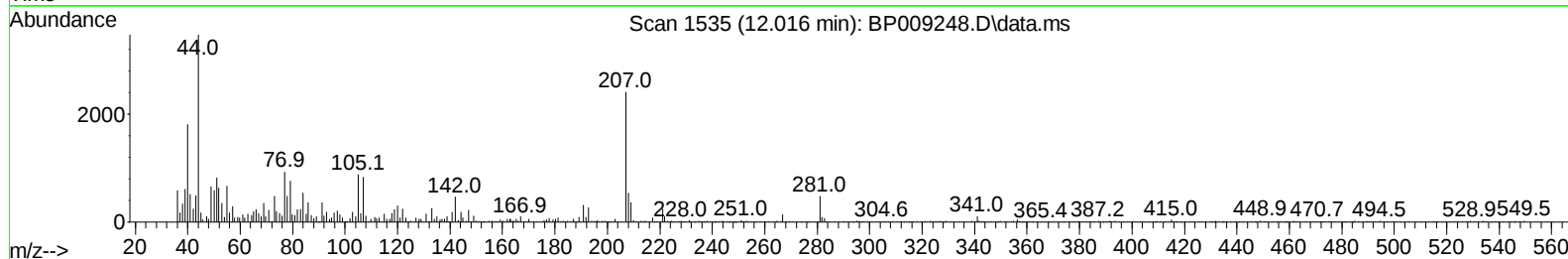
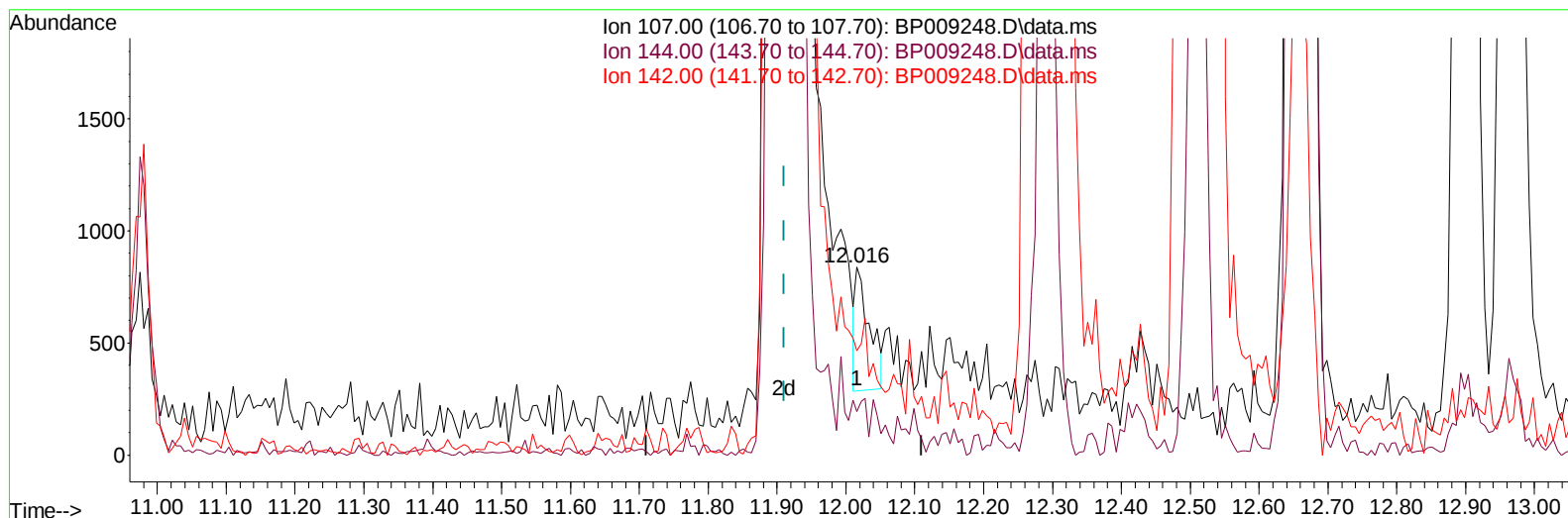
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP030222\  
Data File : BP009248.D  
Acq On : 02 Mar 2022 16:39  
Operator : CG/JU  
Sample : SST02010  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SST020614

Manual IntegrationsAPPROVED

Quant Time: Mar 02 17:37:09 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030222.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Mar 02 17:32:26 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/03/2022  
Supervised By :mohammad ahmed 03/07/2022



TIC: BP009248.D\data.ms

(35) 4-Chloro-3-methylphenol (C)

12.016min (+ 0.106) 0.04 ng/u1

response 800

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 107.00 | 100.00 | 100.00 |
| 144.00 | 19.80  | 23.00  |
| 142.00 | 48.60  | 55.54  |
| 0.00   | 0.00   | 0.00   |

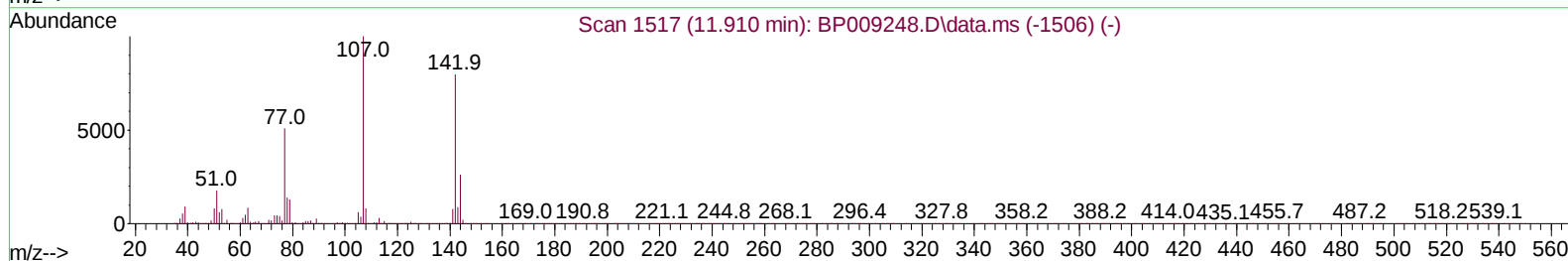
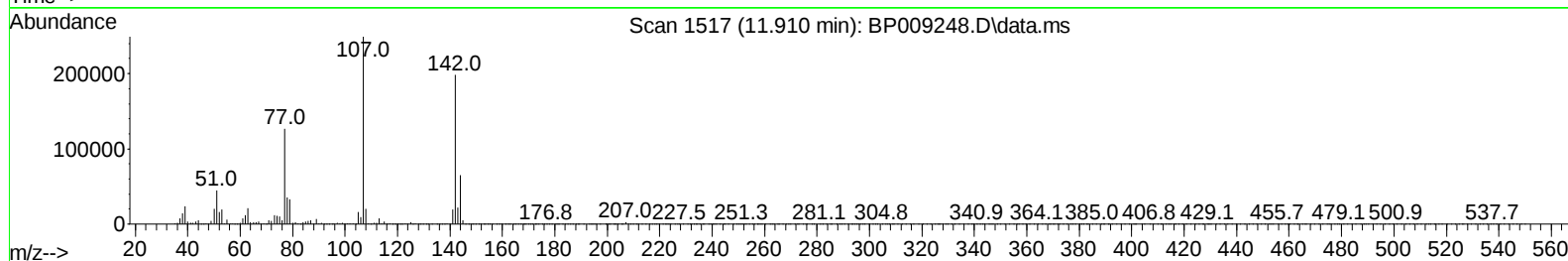
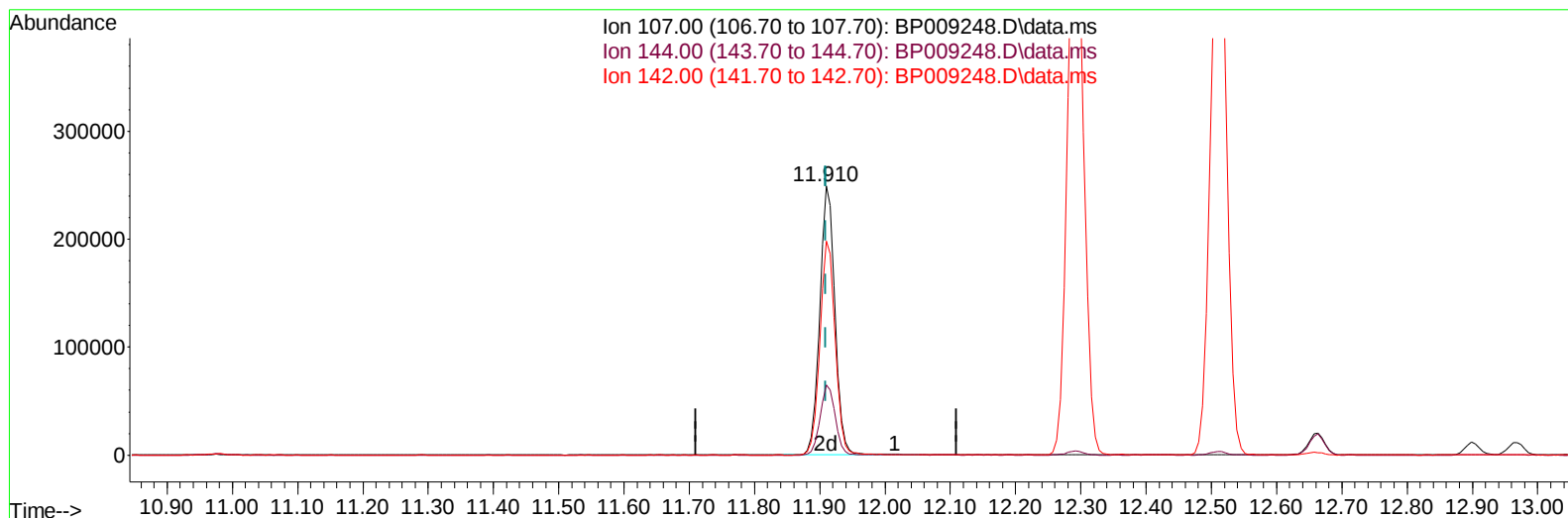
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP030222\  
 Data File : BP009248.D  
 Acq On : 02 Mar 2022 16:39  
 Operator : CG/JU  
 Sample : SSTD02010  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD020614

Manual IntegrationsAPPROVED

Quant Time: Mar 02 17:37:09 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030222.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Mar 02 17:32:26 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/03/2022  
 Supervised By :mohammad ahmed 03/07/2022



TIC: BP009248.D\data.ms

(35) 4-Chloro-3-methylphenol (C)

11.910min ( 0.000) 20.50 ng/ul m

response 407067

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 107.00 | 100.00 | 100.00 |
| 144.00 | 19.80  | 26.00# |
| 142.00 | 48.60  | 79.60# |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP030222\  
 Data File : BP009248.D  
 Acq On : 02 Mar 2022 16:39  
 Operator : CG/JU  
 Sample : SST002010  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST0020614

Manual IntegrationsAPPROVED

Quant Time: Mar 02 17:37:09 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030222.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Mar 02 17:32:26 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/03/2022  
 Supervised By :mohammad ahmed 03/07/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.828  | 152  | 349859   | 20.000 | ng/ul  | # 0.00   |
| 20) Naphthalene-d8            | 10.628 | 136  | 1357567  | 20.000 | ng/ul  | # 0.00   |
| 38) Acenaphthene-d10          | 14.475 | 164  | 791554   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.233 | 188  | 1553347  | 20.000 | ng/ul  | # 0.00   |
| 79) Chrysene-d12              | 21.339 | 240  | 1553653  | 20.000 | ng/ul  | # 0.00   |
| 88) Perylene-d12              | 23.756 | 264  | 1674106  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.322  | 96   | 71704m   | 8.398  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.722  | 84   | 430087   | 19.842 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.993  | 99   | 556179   | 19.654 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.163  | 67   | 333688   | 19.570 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.363  | 132  | 441596   | 19.913 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.534  | 113  | 424412   | 19.532 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.981  | 128  | 172526   | 22.423 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.704  | 143  | 167791   | 23.264 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.245 | 165  | 411805   | 20.690 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.757 | 131  | 568856   | 20.142 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.886 | 166  | 1120465  | 20.212 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.163 | 160  | 1465999  | 19.957 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.657 | 143  | 176283   | 20.861 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.469 | 176  | 971027   | 20.193 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.580 | 200  | 116672   | 19.341 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.333 | 188  | 1473143  | 20.202 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.574 | 212  | 1698737  | 19.777 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.598 | 264  | 1664196  | 19.616 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.352  | 88   | 68769    | 7.938  | ng/ul# | 49       |
| 5) Pyridine                   | 3.740  | 79   | 462033   | 20.177 | ng/ul# | 75       |
| 6) Benzaldehyde               | 6.969  | 77   | 383958m  | 22.909 | ng/ul  |          |
| 8) Phenol                     | 7.022  | 94   | 563044   | 19.750 | ng/ul  | 90       |
| 10) Bis(2-Chloroethyl)ether   | 7.257  | 93   | 445854   | 19.517 | ng/ul# | 86       |
| 12) 2-Chlorophenol            | 7.393  | 128  | 453227   | 20.225 | ng/ul# | 88       |
| 13) 2-Methylphenol            | 8.269  | 108  | 411592   | 19.205 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.363  | 45   | 540033   | 19.591 | ng/ul# | 91       |
| 16) Acetophenone              | 8.651  | 105  | 658607   | 19.704 | ng/ul# | 87       |
| 17) N-Nitroso-di-n-propyla... | 8.640  | 70   | 326829   | 19.603 | ng/ul# | 83       |
| 18) 4-Methylphenol            | 8.598  | 108  | 443375   | 19.718 | ng/ul  | 92       |
| 19) Hexachloroethane          | 8.910  | 117  | 181343   | 19.770 | ng/ul# | 77       |
| 22) Nitrobenzene              | 9.022  | 77   | 451571   | 22.243 | ng/ul# | 85       |
| 23) Isophorone                | 9.557  | 82   | 902096   | 19.640 | ng/ul# | 95       |
| 25) 2-Nitrophenol             | 9.734  | 139  | 197969   | 23.177 | ng/ul# | 75       |
| 26) 2,4-Dimethylphenol        | 9.804  | 107  | 479174   | 20.211 | ng/ul# | 71       |
| 27) Bis(2-Chloroethoxy)met... | 10.040 | 93   | 558198   | 19.952 | ng/ul# | 97       |
| 29) 2,4-Dichlorophenol        | 10.269 | 162  | 403569   | 20.551 | ng/ul# | 83       |
| 30) Naphthalene               | 10.675 | 128  | 1377254  | 19.982 | ng/ul  | 96       |
| 32) 4-Chloroaniline           | 10.781 | 127  | 569715   | 20.435 | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 10.981 | 225  | 282386   | 19.809 | ng/ul  | 95       |
| 34) Caprolactam               | 11.545 | 113  | 108461   | 19.133 | ng/ul# | 46       |
| 35) 4-Chloro-3-methylphenol   | 11.910 | 107  | 407067m  | 20.502 | ng/ul  |          |

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 Sample : SST02010  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST020614

## Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 03/07/2022

Quant Time: Mar 02 17:37:09 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030222.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Mar 02 17:32:26 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.292 | 142  | 917869   | 19.818 | ng/ul  | 91       |
| 37) 1-Methylnaphthalene       | 12.510 | 142  | 929478   | 20.088 | ng/ul# | 96       |
| 39) 1,2,4,5-Tetrachloroben... | 12.663 | 216  | 501536   | 19.905 | ng/ul  | 94       |
| 40) Hexachlorocyclopentadiene | 12.651 | 237  | 303738   | 19.365 | ng/ul  | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.898 | 196  | 300296   | 20.628 | ng/ul# | 83       |
| 42) 2,4,5-Trichlorophenol     | 12.969 | 196  | 316465   | 20.544 | ng/ul# | 86       |
| 43) 1,1'-Biphenyl             | 13.310 | 154  | 1201463  | 19.836 | ng/ul# | 94       |
| 44) 2-Chloronaphthalene       | 13.345 | 162  | 920098   | 19.982 | ng/ul  | 92       |
| 45) 2-Nitroaniline            | 13.539 | 65   | 176338   | 22.640 | ng/ul# | 85       |
| 47) Dimethylphthalate         | 13.933 | 163  | 1086929  | 20.309 | ng/ul# | 92       |
| 48) 2,6-Dinitrotoluene        | 14.039 | 165  | 162350   | 22.826 | ng/ul  | 97       |
| 50) Acenaphthylene            | 14.192 | 152  | 1471574  | 20.270 | ng/ul  | 95       |
| 51) 3-Nitroaniline            | 14.369 | 138  | 188033   | 20.981 | ng/ul  | 80       |
| 52) Acenaphthene              | 14.533 | 153  | 940411   | 19.915 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.575 | 184  | 120607   | 31.880 | ng/ul# | 82       |
| 55) 4-Nitrophenol             | 14.675 | 109  | 139122   | 20.591 | ng/ul# | 34       |
| 56) Dibenzofuran              | 14.875 | 168  | 1343838  | 20.146 | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene        | 14.828 | 165  | 243243   | 24.073 | ng/ul  | 91       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.098 | 232  | 254492   | 21.347 | ng/ul# | 86       |
| 59) Diethylphthalate          | 15.304 | 149  | 1056971  | 20.438 | ng/ul  | 93       |
| 61) Fluorene                  | 15.527 | 166  | 1065665  | 20.436 | ng/ul  | 90       |
| 62) 4-Chlorophenyl-phenyle... | 15.527 | 204  | 535513   | 20.220 | ng/ul  | 97       |
| 63) 4-Nitroaniline            | 15.533 | 138  | 195721   | 22.664 | ng/ul# | 71       |
| 66) 4,6-Dinitro-2-methylph... | 15.598 | 198  | 127031   | 20.072 | ng/ul# | 94       |
| 67) N-Nitrosodiphenylamine    | 15.733 | 169  | 919059   | 20.090 | ng/ul  | 94       |
| 68) 4-Bromophenyl-phenylether | 16.422 | 248  | 321835   | 19.722 | ng/ul  | 94       |
| 69) Hexachlorobenzene         | 16.539 | 284  | 368591   | 19.768 | ng/ul# | 89       |
| 70) Atrazine                  | 16.698 | 200  | 319133   | 19.696 | ng/ul# | 90       |
| 71) Pentachlorophenol         | 16.880 | 266  | 207058   | 18.976 | ng/ul# | 85       |
| 72) Phenanthrene              | 17.274 | 178  | 1660516  | 20.083 | ng/ul  | 98       |
| 74) Anthracene                | 17.369 | 178  | 1674069  | 20.288 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.269 | 216  | 510172   | 19.183 | ng/uL# | 87       |
| 76) Pentachlorobenzene        | 14.798 | 250  | 452698   | 19.176 | ng/uL  | 88       |
| 77) Carbazole                 | 17.639 | 167  | 1510366  | 20.629 | ng/ul# | 96       |
| 78) Di-n-butylphthalate       | 18.210 | 149  | 1696015  | 20.797 | ng/ul# | 92       |
| 80) Fluoranthene              | 19.251 | 202  | 1960632  | 19.793 | ng/ul  | 98       |
| 82) Pyrene                    | 19.604 | 202  | 1990821  | 19.695 | ng/ul  | 98       |
| 83) Butylbenzylphthalate      | 20.498 | 149  | 694101   | 21.538 | ng/ul# | 80       |
| 84) 3,3'-Dichlorobenzidine    | 21.257 | 252  | 665204   | 20.599 | ng/ul# | 96       |
| 85) Benzo(a)anthracene        | 21.327 | 228  | 1966064  | 20.033 | ng/ul  | 95       |
| 86) Bis(2-ethylhexyl)phtha... | 21.274 | 149  | 1089987  | 21.147 | ng/ul# | 80       |
| 87) Chrysene                  | 21.380 | 228  | 1905850  | 20.058 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.204 | 149  | 1735059  | 19.176 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.021 | 252  | 2068435  | 19.577 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.068 | 252  | 1918275  | 20.137 | ng/ul# | 96       |
| 93) Benzo(a)pyrene            | 23.651 | 252  | 2003944  | 19.820 | ng/ul# | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.268 | 276  | 2353257  | 19.686 | ng/ul# | 92       |
| 95) Dibenzo(a,h)anthracene    | 26.292 | 278  | 2002060  | 19.823 | ng/ul# | 92       |
| 96) Benzo(g,h,i)perylene      | 27.039 | 276  | 1998891  | 19.478 | ng/ul# | 86       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SSTD020614

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 03/03/2022  
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