

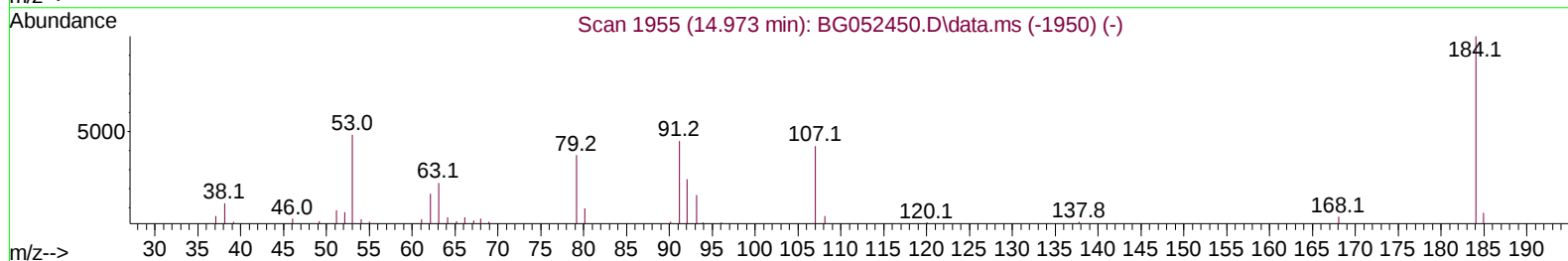
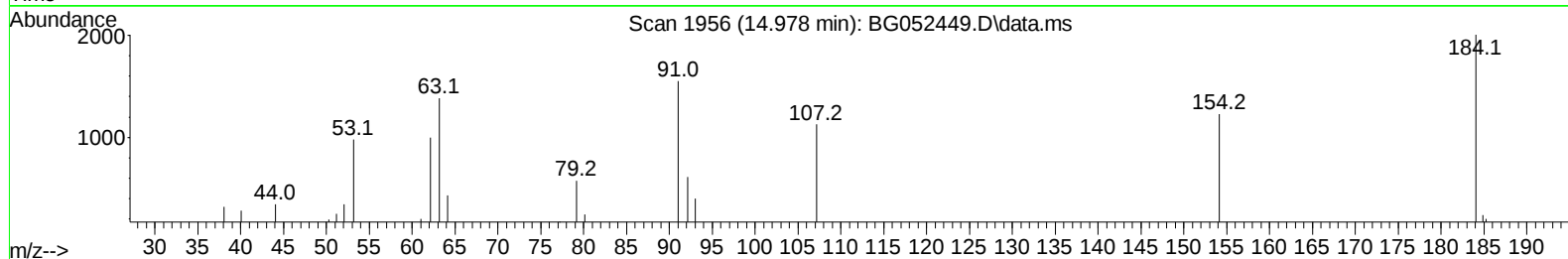
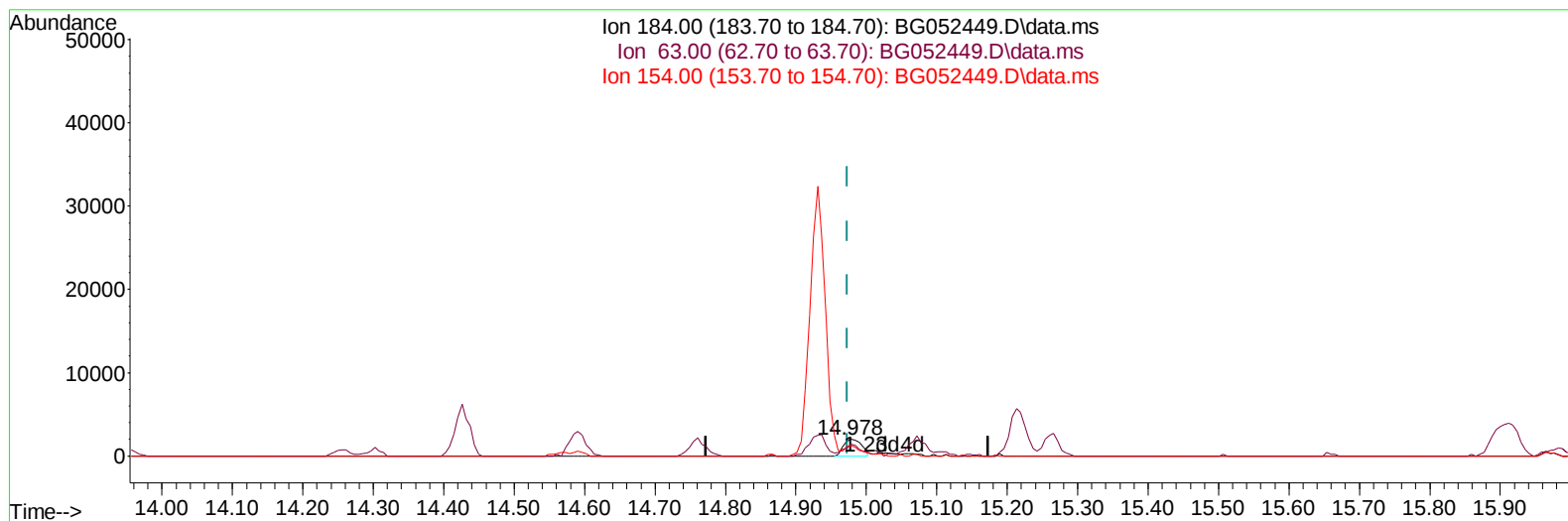
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022222\  
Data File : BG052449.D  
Acq On : 22 Feb 2022 14:06  
Operator : CG/JU  
Sample : SST01008  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SST010408

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/23/2022  
Supervised By :Yogesh Patel 02/24/2022

Quant Time: Feb 22 15:25:23 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022222.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Feb 22 15:23:39 2022  
Response via : Initial Calibration



TIC: BG052449.D\data.ms

**(53) 2,4-Dinitrophenol****14.978min (+ 0.005) 5.40 ng/ul****response 3561**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 184.00 | 100.00 | 100.00 |
| 63.00  | 82.70  | 68.92  |
| 154.00 | 67.00  | 61.21  |
| 0.00   | 0.00   | 0.00   |

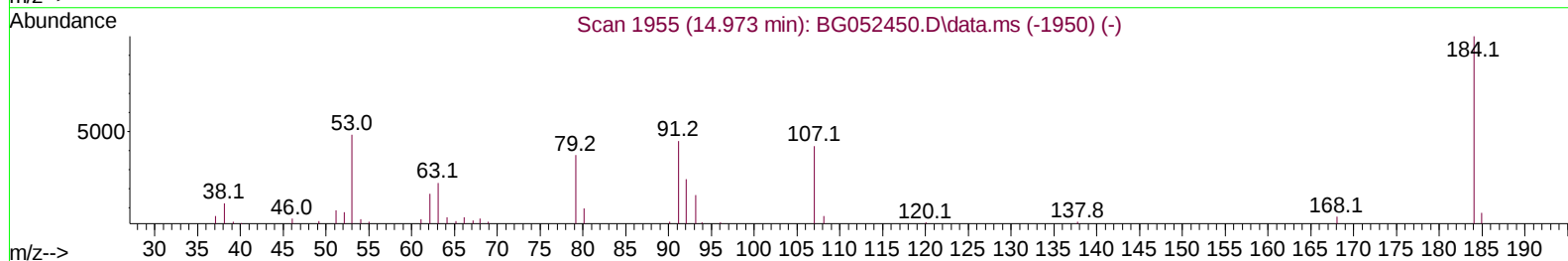
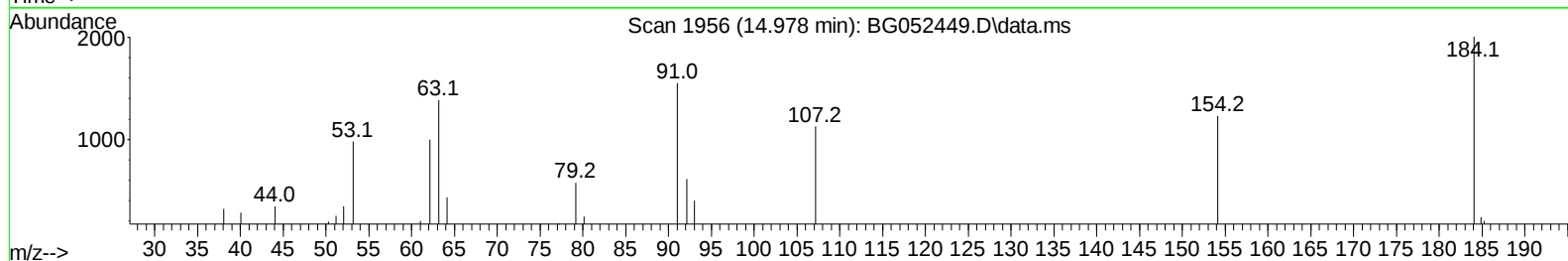
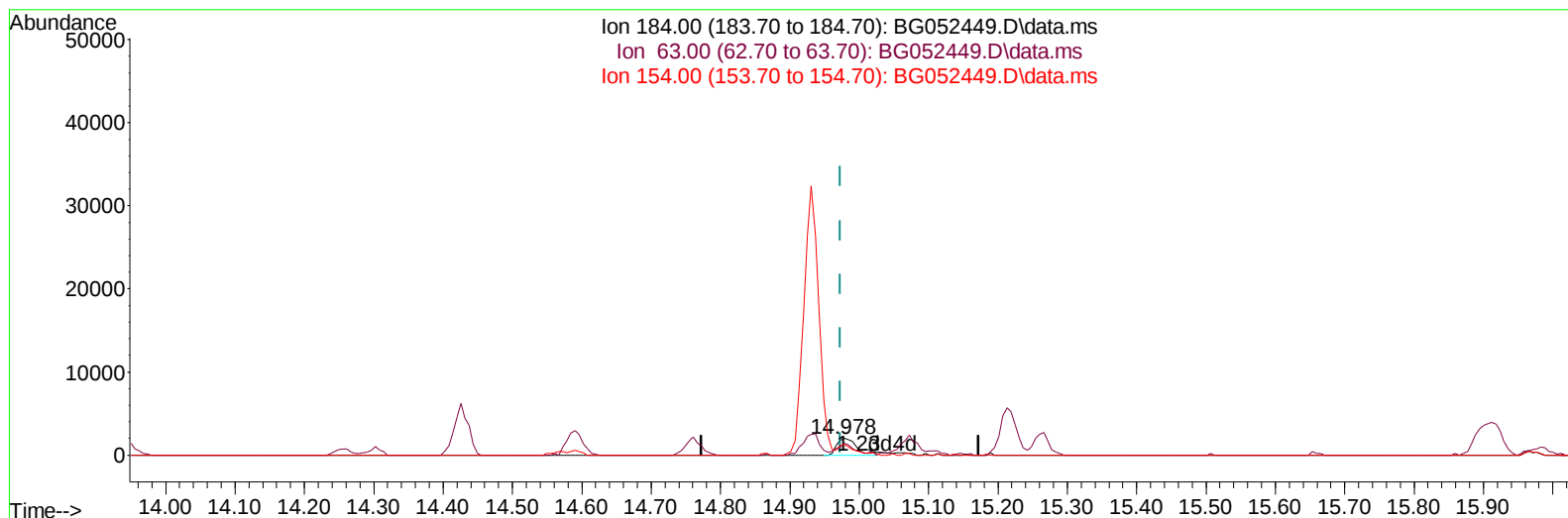
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022222\  
 Data File : BG052449.D  
 Acq On : 22 Feb 2022 14:06  
 Operator : CG/JU  
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 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST010408

Manual IntegrationsAPPROVED

Quant Time: Feb 22 15:25:23 2022  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 22 15:23:39 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/23/2022  
 Supervised By :Yogesh Patel 02/24/2022



TIC: BG052449.D\data.ms

**(53) 2,4-Dinitrophenol**

14.978min (+ 0.005) 6.42 ng/ul m

response 4236

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 184.00 | 100.00 | 100.00 |
| 63.00  | 82.70  | 68.92  |
| 154.00 | 67.00  | 61.21  |
| 0.00   | 0.00   | 0.00   |

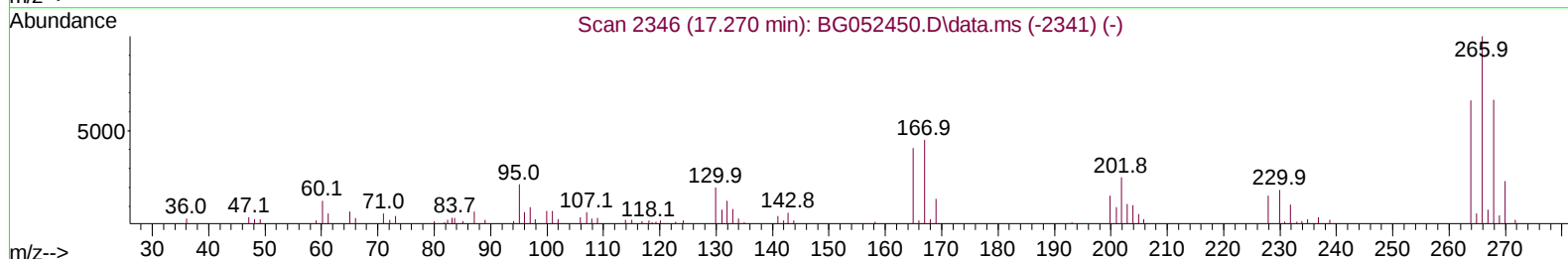
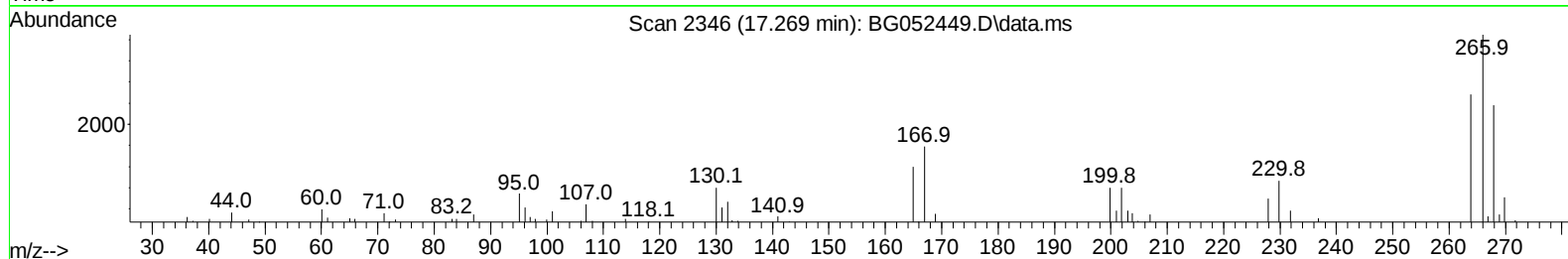
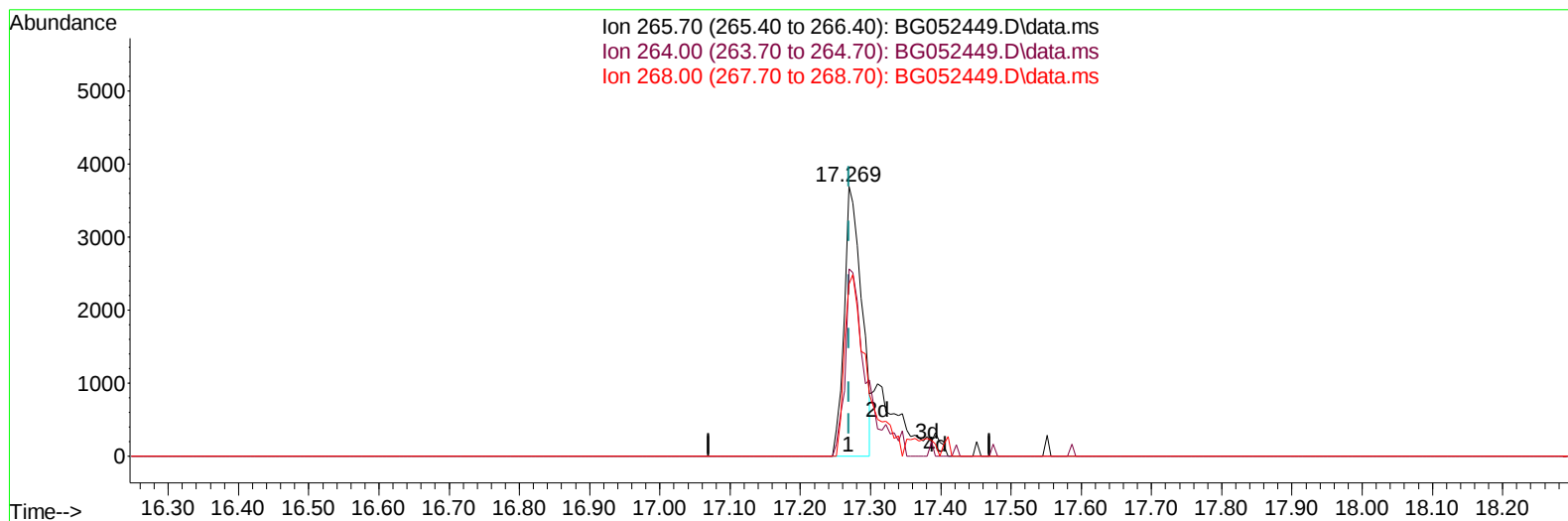
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022222\  
 Data File : BG052449.D  
 Acq On : 22 Feb 2022 14:06  
 Operator : CG/JU  
 Sample : SST01008  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST010408

Manual IntegrationsAPPROVED

Quant Time: Feb 22 15:25:23 2022  
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 Supervised By :Yogesh Patel 02/24/2022



TIC: BG052449.D\data.ms

**(71) Pentachlorophenol (C)**

**17.269min (-0.001) 5.01 ng/u1**

**response 6323**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 265.70 | 100.00 | 100.00 |
| 264.00 | 67.90  | 69.54  |
| 268.00 | 63.80  | 63.90  |
| 0.00   | 0.00   | 0.00   |

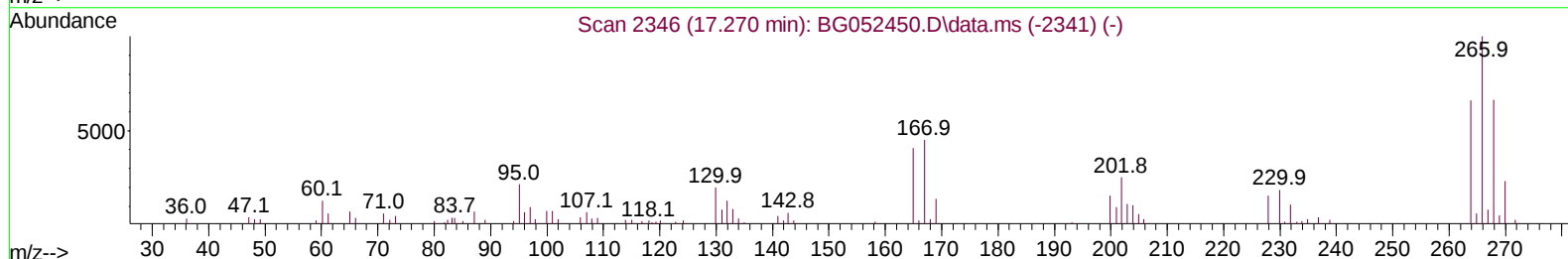
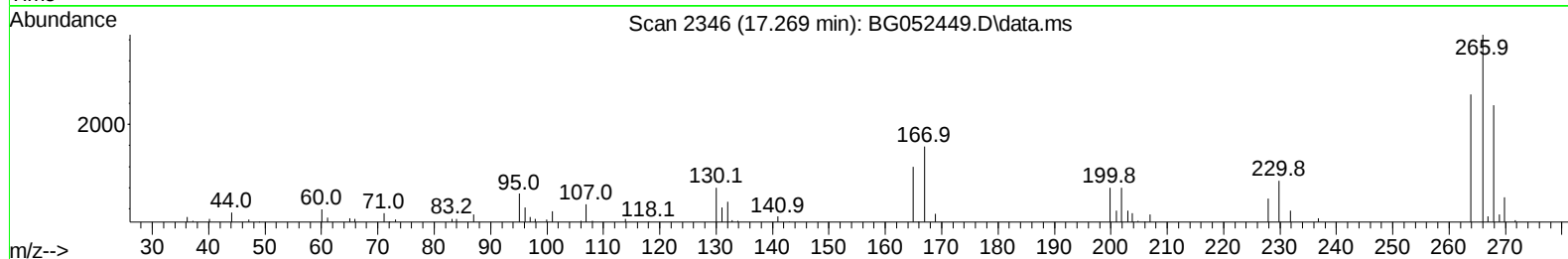
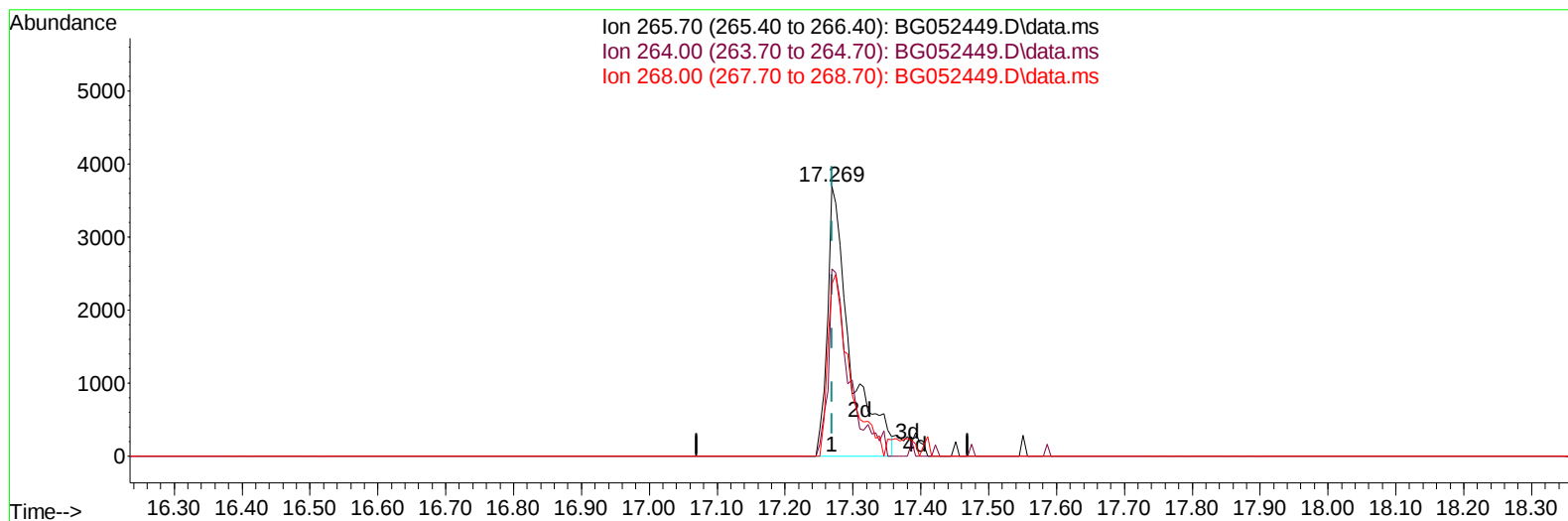
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022222\  
Data File : BG052449.D  
Acq On : 22 Feb 2022 14:06  
Operator : CG/JU  
Sample : SST01008  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD010408

Manual IntegrationsAPPROVED

Quant Time: Feb 22 15:25:23 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022222.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Feb 22 15:23:39 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/23/2022  
Supervised By :Yogesh Patel 02/24/2022



TIC: BG052449.D\data.ms

**(71) Pentachlorophenol (C)**

**17.269min (-0.001) 6.78 ng/ul m**

**response 8551**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 265.70 | 100.00 | 100.00 |
| 264.00 | 67.90  | 69.54  |
| 268.00 | 63.80  | 63.90  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022222\  
 Data File : BG052449.D  
 Acq On : 22 Feb 2022 14:06  
 Operator : CG/JU  
 Sample : SST001008  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

SSTD010408

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/23/2022  
 Supervised By :Yogesh Patel 02/24/2022

Quant Time: Feb 22 15:25:23 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022222.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 22 15:23:39 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.222  | 152  | 26287    | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 11.060 | 136  | 110953   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.866 | 164  | 77277    | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.616 | 188  | 193876   | 20.000 | ng/ul  | # 0.00   |
| 79) Chrysene-d12              | 21.933 | 240  | 203405   | 20.000 | ng/ul  | # 0.00   |
| 88) Perylene-d12              | 25.393 | 264  | 204528   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.546  | 96   | 2972     | 4.561  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.981  | 84   | 20142    | 9.938  | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.371  | 99   | 26152    | 10.937 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.529  | 67   | 15470    | 10.650 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.752  | 132  | 18801    | 11.284 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.927  | 113  | 19574    | 10.788 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.391  | 128  | 10410    | 11.523 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.126 | 143  | 10794    | 10.934 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.678 | 165  | 20213    | 10.503 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.189 | 131  | 27385    | 11.218 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.255 | 166  | 68888    | 11.341 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.561 | 160  | 84410    | 11.149 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.054 | 143  | 9359     | 10.173 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.853 | 176  | 58976    | 11.082 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.971 | 200  | 11211    | 9.366  | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.715 | 188  | 102642   | 11.230 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.995 | 212  | 128053   | 10.504 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.152 | 264  | 117836   | 10.845 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.588  | 88   | 3523     | 4.851  | ng/ul# | 90       |
| 5) Pyridine                   | 3.999  | 79   | 20806    | 9.803  | ng/ul# | 84       |
| 6) Benzaldehyde               | 7.341  | 77   | 18833    | 13.893 | ng/ul# | 89       |
| 8) Phenol                     | 7.394  | 94   | 26970    | 11.255 | ng/ul  | 93       |
| 10) Bis(2-Chloroethyl)ether   | 7.623  | 93   | 19812    | 10.941 | ng/ul  | 95       |
| 12) 2-Chlorophenol            | 7.782  | 128  | 19399    | 11.279 | ng/ul  | 93       |
| 13) 2-Methylphenol            | 8.663  | 108  | 19461    | 10.791 | ng/ul  | 97       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.745  | 45   | 28162    | 10.889 | ng/ul# | 95       |
| 16) Acetophenone              | 9.045  | 105  | 31804    | 11.303 | ng/ul  | 94       |
| 17) N-Nitroso-di-n-propyla... | 9.021  | 70   | 19494    | 11.549 | ng/ul  | 96       |
| 18) 4-Methylphenol            | 8.992  | 108  | 20703    | 10.897 | ng/ul  | 94       |
| 19) Hexachloroethane          | 9.321  | 117  | 7709     | 10.407 | ng/ul# | 81       |
| 22) Nitrobenzene              | 9.438  | 77   | 26251    | 10.418 | ng/ul  | 96       |
| 23) Isophorone                | 9.961  | 82   | 49700    | 10.659 | ng/ul# | 99       |
| 25) 2-Nitrophenol             | 10.161 | 139  | 11674    | 10.817 | ng/ul  | 93       |
| 26) 2,4-Dimethylphenol        | 10.208 | 107  | 24328    | 10.932 | ng/ul  | 94       |
| 27) Bis(2-Chloroethoxy)met... | 10.437 | 93   | 27312    | 10.922 | ng/ul  | 97       |
| 29) 2,4-Dichlorophenol        | 10.707 | 162  | 19398    | 10.427 | ng/ul  | 94       |
| 30) Naphthalene               | 11.113 | 128  | 67681    | 11.156 | ng/ul# | 94       |
| 32) 4-Chloroaniline           | 11.213 | 127  | 28473    | 11.313 | ng/ul  | 94       |
| 33) Hexachlorobutadiene       | 11.400 | 225  | 15954    | 10.513 | ng/ul  | 98       |
| 34) Caprolactam               | 11.953 | 113  | 6771     | 11.222 | ng/ul  | 86       |
| 35) 4-Chloro-3-methylphenol   | 12.329 | 107  | 20896    | 10.414 | ng/ul  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG022222\  
 Data File : BG052449.D  
 Acq On : 22 Feb 2022 14:06  
 Operator : CG/JU  
 Sample : SST001008  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

SSTD010408

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/23/2022  
 Supervised By :Yogesh Patel 02/24/2022

Quant Time: Feb 22 15:25:23 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022222.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 22 15:23:39 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.705 | 142  | 48307    | 11.347 | ng/ul  | 90       |
| 37) 1-Methylnaphthalene       | 12.928 | 142  | 48969    | 11.181 | ng/ul  | 97       |
| 39) 1,2,4,5-Tetrachloroben... | 13.075 | 216  | 30025    | 10.794 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 13.051 | 237  | 10083    | 6.350  | ng/ul# | 92       |
| 41) 2,4,6-Trichlorophenol     | 13.310 | 196  | 17483    | 10.527 | ng/ul  | 98       |
| 42) 2,4,5-Trichlorophenol     | 13.386 | 196  | 18156    | 10.150 | ng/ul  | 96       |
| 43) 1,1'-Biphenyl             | 13.697 | 154  | 67515    | 11.144 | ng/ul# | 94       |
| 44) 2-Chloronaphthalene       | 13.750 | 162  | 51428    | 11.095 | ng/ul  | 96       |
| 45) 2-Nitroaniline            | 13.944 | 65   | 15938    | 10.212 | ng/ul# | 88       |
| 47) Dimethylphthalate         | 14.302 | 163  | 66885    | 11.360 | ng/ul# | 98       |
| 48) 2,6-Dinitrotoluene        | 14.426 | 165  | 14962    | 11.780 | ng/ul# | 91       |
| 50) Acenaphthylene            | 14.590 | 152  | 85949    | 11.326 | ng/ul  | 96       |
| 51) 3-Nitroaniline            | 14.761 | 138  | 13336    | 12.269 | ng/ul  | 96       |
| 52) Acenaphthene              | 14.931 | 153  | 55986    | 11.254 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.978 | 184  | 4236m    | 6.422  | ng/ul  |          |
| 55) 4-Nitrophenol             | 15.072 | 109  | 9510     | 9.863  | ng/ul  | 86       |
| 56) Dibenzofuran              | 15.266 | 168  | 80348    | 11.155 | ng/ul  | 100      |
| 57) 2,4-Dinitrotoluene        | 15.213 | 165  | 20842    | 11.447 | ng/ul# | 84       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.495 | 232  | 16579    | 10.588 | ng/ul# | 86       |
| 59) Diethylphthalate          | 15.659 | 149  | 67571    | 11.318 | ng/ul  | 99       |
| 61) Fluorene                  | 15.912 | 166  | 68556    | 11.469 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.894 | 204  | 37358    | 11.364 | ng/ul  | 89       |
| 63) 4-Nitroaniline            | 15.918 | 138  | 14481    | 13.730 | ng/ul  | 96       |
| 66) 4,6-Dinitro-2-methylph... | 15.988 | 198  | 10979    | 9.631  | ng/ul# | 84       |
| 67) N-Nitrosodiphenylamine    | 16.106 | 169  | 59808    | 11.252 | ng/ul  | 93       |
| 68) 4-Bromophenyl-phenylether | 16.793 | 248  | 24357    | 10.407 | ng/ul# | 86       |
| 69) Hexachlorobenzene         | 16.922 | 284  | 27937    | 10.621 | ng/ul# | 88       |
| 70) Atrazine                  | 17.046 | 200  | 26659    | 11.369 | ng/ul  | 96       |
| 71) Pentachlorophenol         | 17.269 | 266  | 8551m    | 6.780  | ng/ul  |          |
| 72) Phenanthrene              | 17.657 | 178  | 113979   | 11.159 | ng/ul  | 97       |
| 74) Anthracene                | 17.751 | 178  | 115126   | 11.420 | ng/ul  | 95       |
| 75) 1,2,3,4-Tetrachloroben... | 13.674 | 216  | 32984    | 10.190 | ng/uL  | 96       |
| 76) Pentachlorobenzene        | 15.189 | 250  | 32273    | 10.808 | ng/uL  | 95       |
| 77) Carbazole                 | 18.015 | 167  | 103633   | 11.795 | ng/ul  | 98       |
| 78) Di-n-butylphthalate       | 18.555 | 149  | 121349   | 11.515 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.660 | 202  | 150239   | 10.465 | ng/ul# | 90       |
| 82) Pyrene                    | 20.024 | 202  | 152265   | 10.834 | ng/ul# | 91       |
| 83) Butylbenzylphthalate      | 20.894 | 149  | 54303    | 11.046 | ng/ul# | 90       |
| 84) 3,3'-Dichlorobenzidine    | 21.810 | 252  | 55850    | 12.823 | ng/ul# | 93       |
| 85) Benzo(a)anthracene        | 21.910 | 228  | 150835   | 11.025 | ng/ul  | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.786 | 149  | 72063    | 10.226 | ng/ul# | 97       |
| 87) Chrysene                  | 21.980 | 228  | 146330   | 11.320 | ng/ul  | 97       |
| 89) Di-n-octyl phthalate      | 23.079 | 149  | 126355   | 10.889 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 24.289 | 252  | 157163   | 11.081 | ng/ul# | 95       |
| 91) Benzo(k)fluoranthene      | 24.359 | 252  | 142255   | 10.963 | ng/ul# | 98       |
| 93) Benzo(a)pyrene            | 25.223 | 252  | 149249   | 11.283 | ng/ul# | 94       |
| 94) Indeno(1,2,3-cd)pyrene    | 29.376 | 276  | 166157   | 11.277 | ng/ul# | 92       |
| 95) Dibenzo(a,h)anthracene    | 29.453 | 278  | 142678   | 11.678 | ng/ul# | 94       |
| 96) Benzo(g,h,i)perylene      | 30.633 | 276  | 138336   | 11.029 | ng/ul# | 85       |

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

**Instrument :**

BNA\_G

**ClientSampleId :**

SSTD010408

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 02/23/2022  
Supervised By :Yogesh Patel 02/24/2022

## Manual IntegrationsAPPROVED

Quant Time: Feb 22 15:25:23 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG022222.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Feb 22 15:23:39 2022  
Response via : Initial Calibration

