

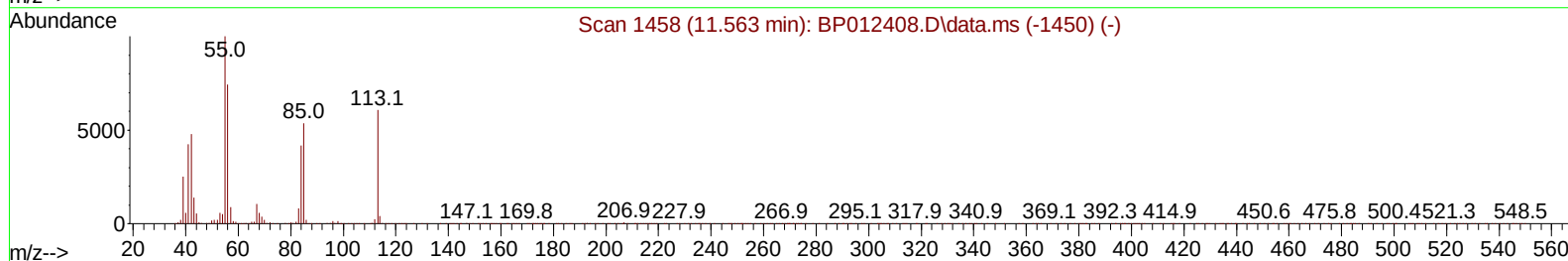
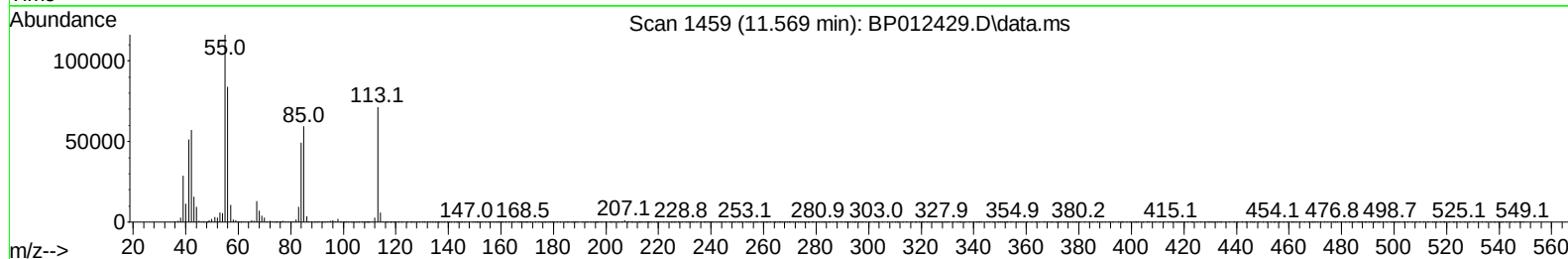
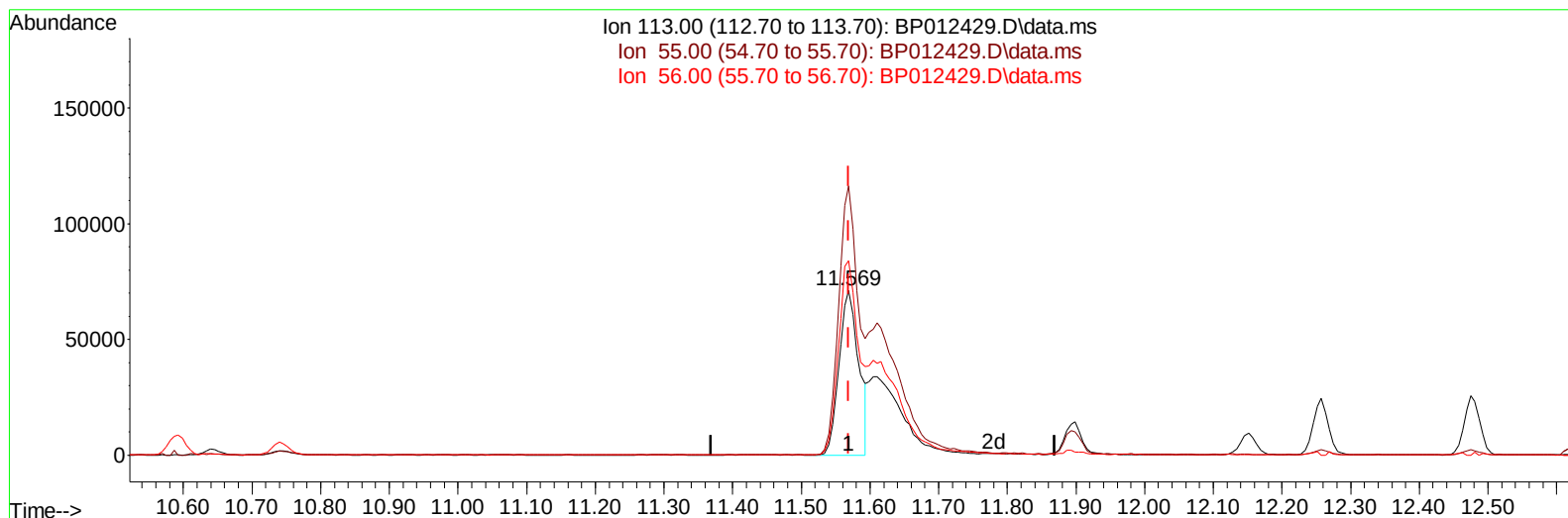
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012429.D  
 Acq On : 01 Nov 2022 16:08  
 Operator : CG/JU  
 Sample : PB148596BS  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS596

Manual IntegrationsAPPROVED

Quant Time: Nov 01 22:59:30 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 01 01:03:58 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/02/2022  
 Supervised By :mohammad ahmed 11/02/2022



TIC: BP012429.D\data.ms

**(34) Caprolactam**

**11.569min (-0.000) 19.33 ng/ul**

**response 141615**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 169.50 | 163.30 |
| 56.00  | 123.60 | 117.87 |
| 0.00   | 0.00   | 0.00   |

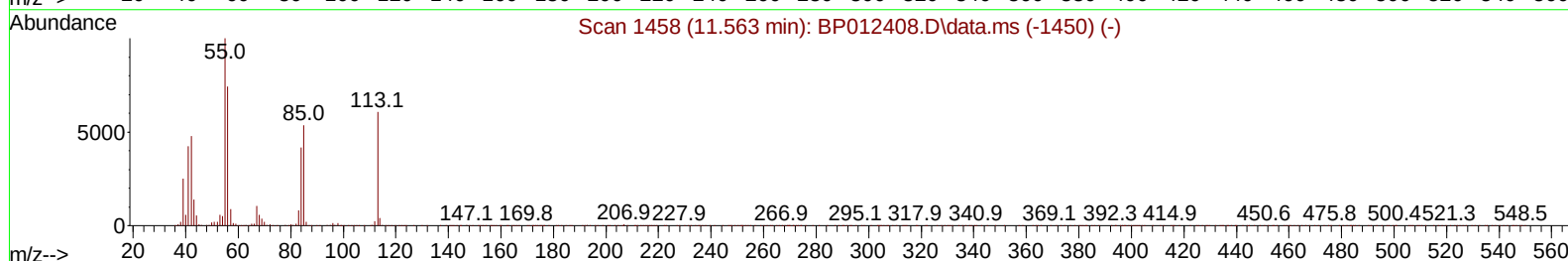
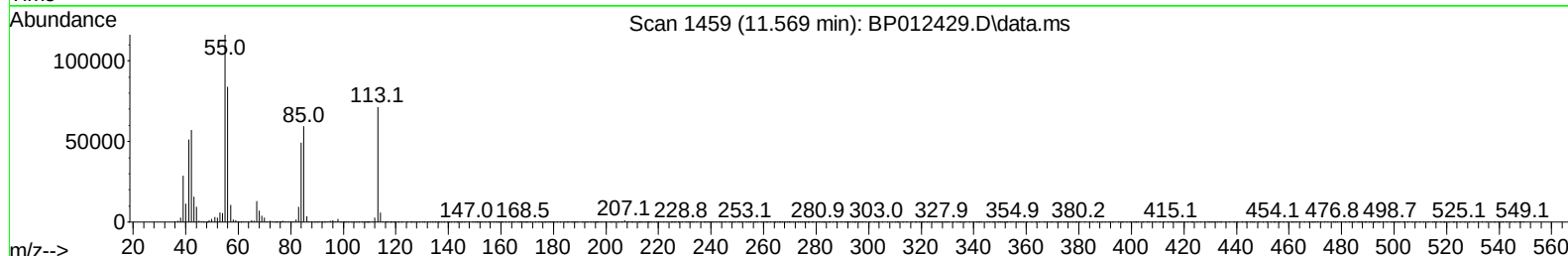
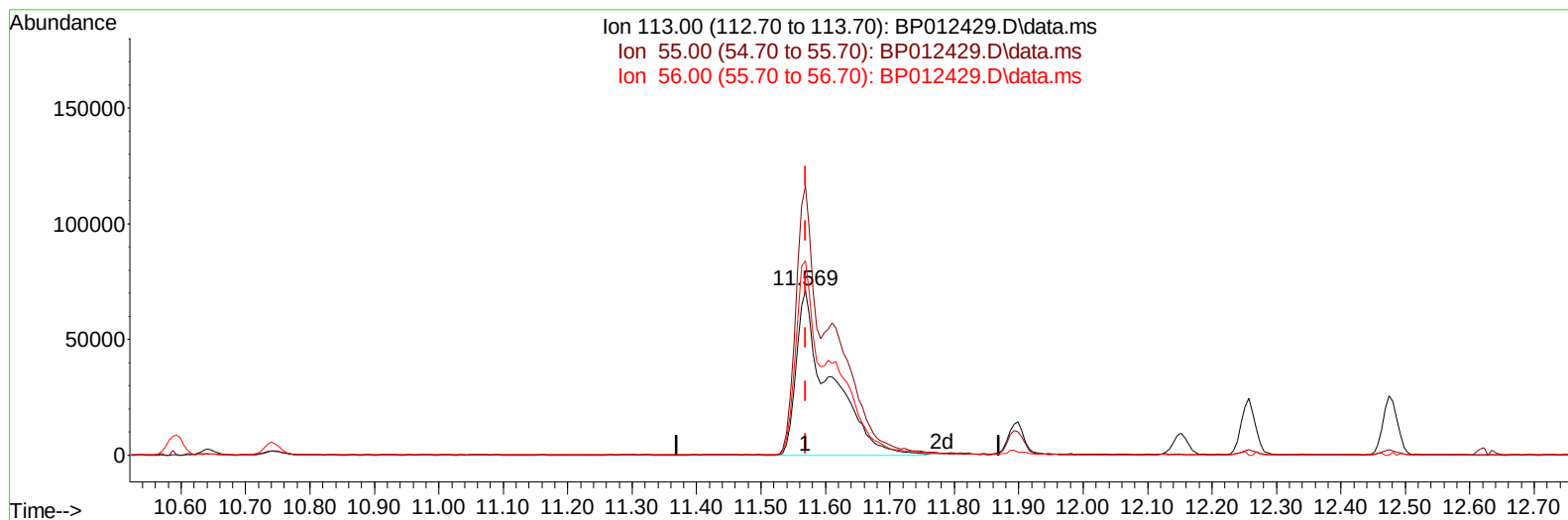
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Instrument :  
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Manual IntegrationsAPPROVED

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TIC: BP012429.D\data.ms

**(34) Caprolactam**

**11.569min (-0.000) 35.30 ng/ul m**

**response 258634**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 169.50 | 163.30 |
| 56.00  | 123.60 | 117.87 |
| 0.00   | 0.00   | 0.00   |

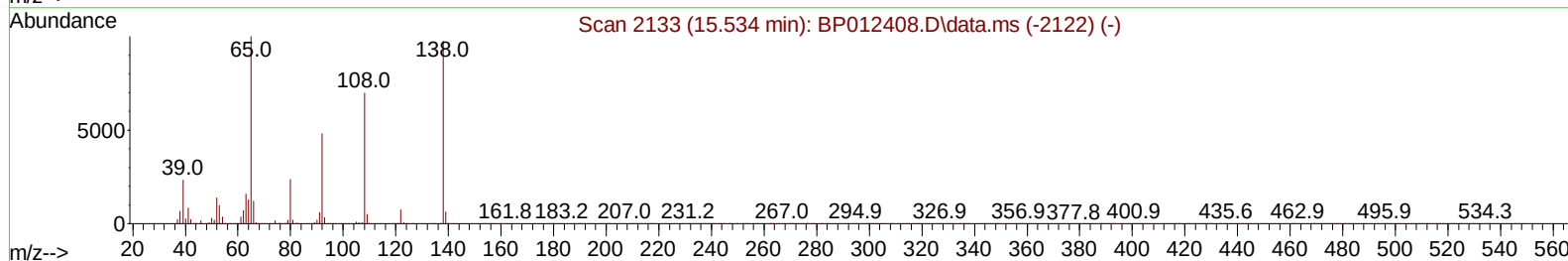
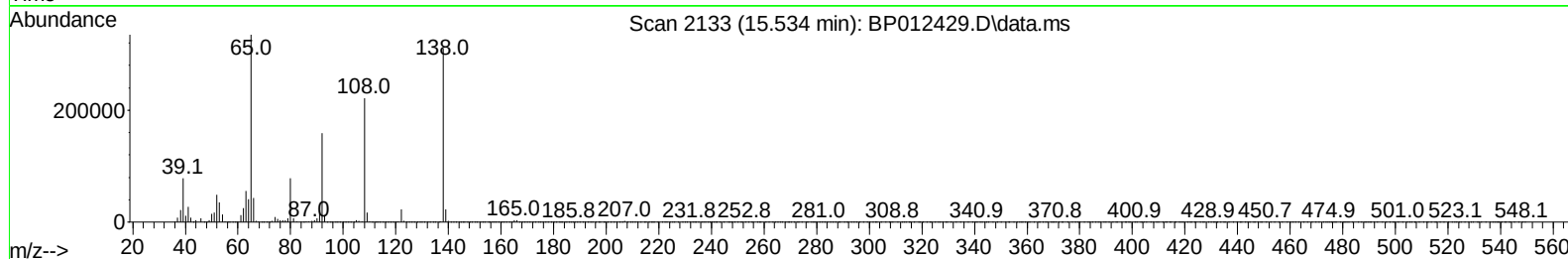
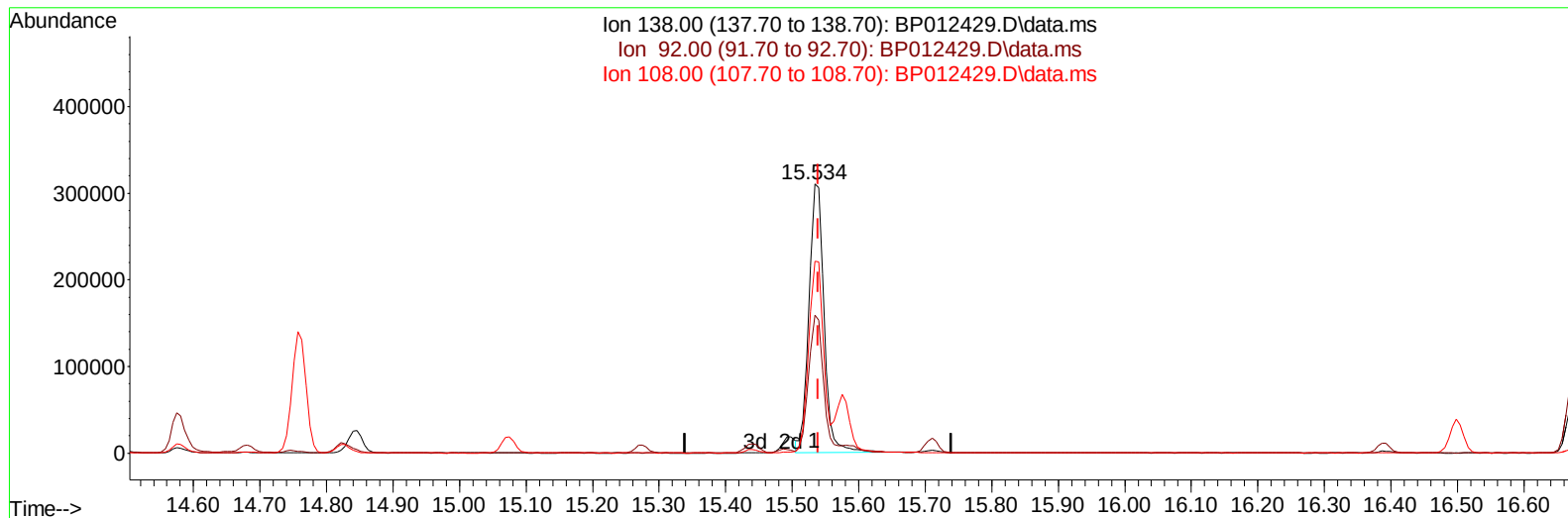
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012429.D  
 Acq On : 01 Nov 2022 16:08  
 Operator : CG/JU  
 Sample : PB148596BS  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
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 ClientSampleId :  
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Manual IntegrationsAPPROVED

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TIC: BP012429.D\data.ms

**(63) 4-Nitroaniline**

**15.534min (-0.006) 39.75 ng/ul**

**response 489782**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 50.80  | 51.30  |
| 108.00 | 70.80  | 71.41  |
| 0.00   | 0.00   | 0.00   |

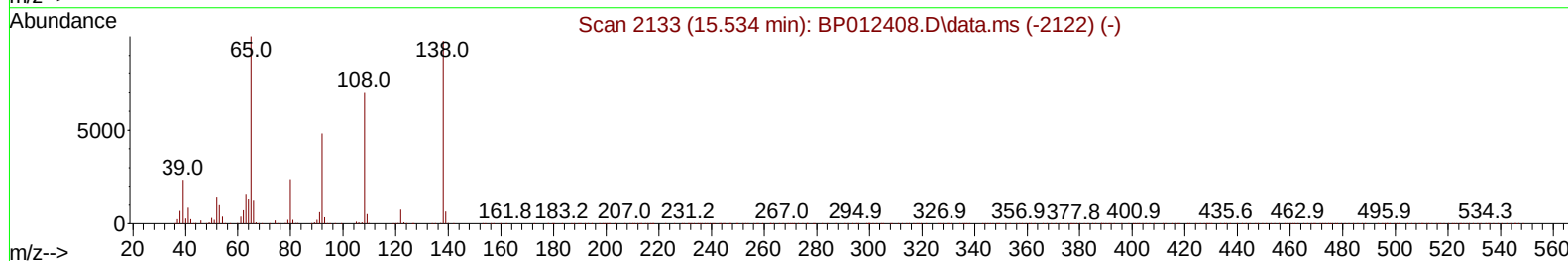
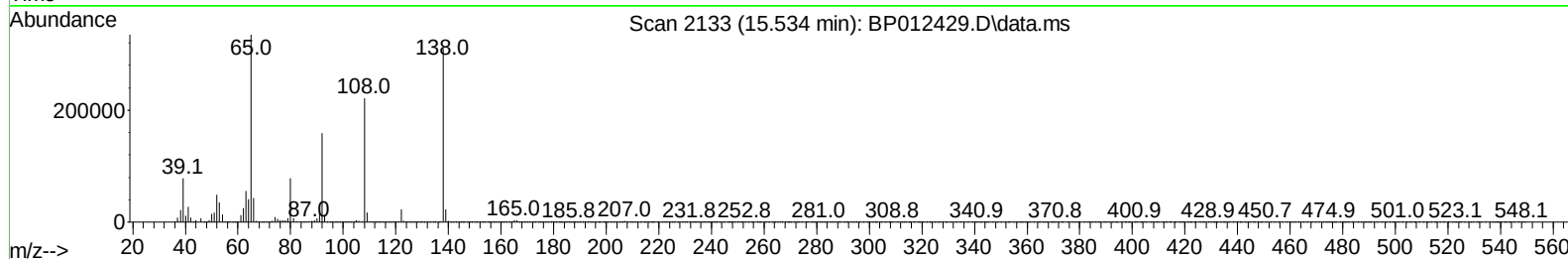
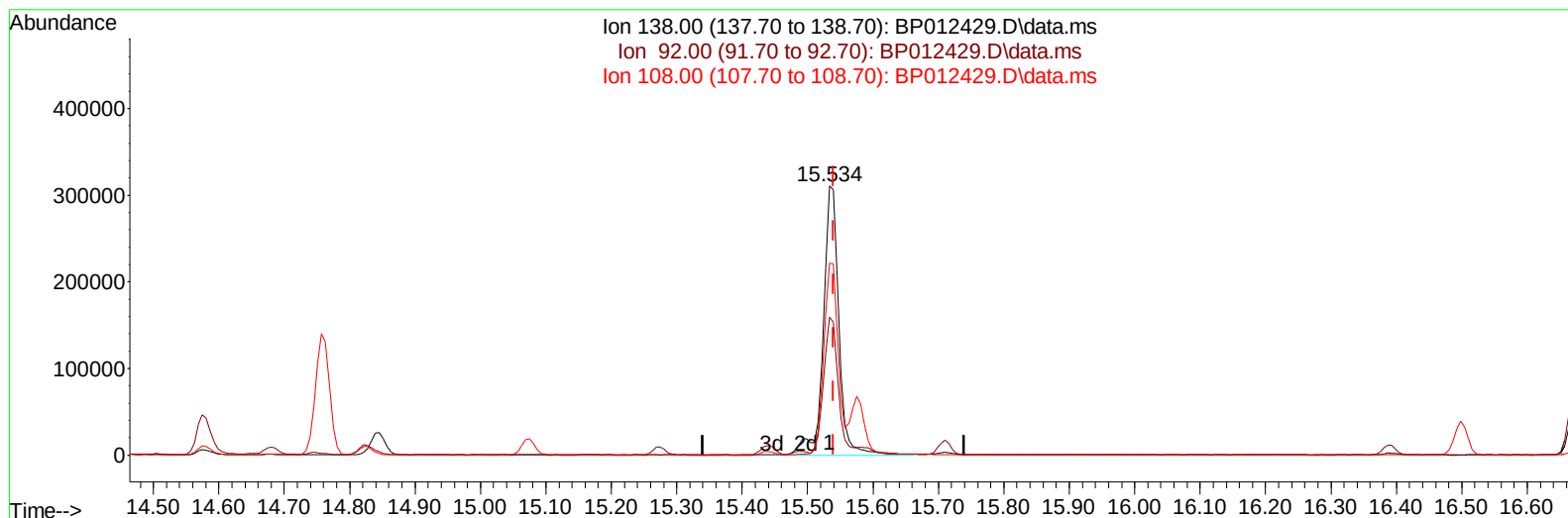
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012429.D  
 Acq On : 01 Nov 2022 16:08  
 Operator : CG/JU  
 Sample : PB148596BS  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
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Manual IntegrationsAPPROVED

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TIC: BP012429.D\data.ms

**(63) 4-Nitroaniline**

**15.534min (-0.006) 42.35 ng/ul m**

**response 521798**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 50.80  | 51.30  |
| 108.00 | 70.80  | 71.41  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012429.D  
 Acq On : 01 Nov 2022 16:08  
 Operator : CG/JU  
 Sample : PB148596BS  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS596

Manual IntegrationsAPPROVED

Quant Time: Nov 01 22:59:30 2022  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 01 01:03:58 2022  
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Reviewed By :Jagrut Upadhyay 11/02/2022  
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| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.793  | 152  | 310354   | 20.000 | ng/uI  | 0.00     |
| 20) Naphthalene-d8            | 10.593 | 136  | 1422167  | 20.000 | ng/uI  | 0.00     |
| 38) Acenaphthene-d10          | 14.440 | 164  | 930064   | 20.000 | ng/uI  | 0.00     |
| 64) Phenanthrene-d10          | 17.204 | 188  | 1990214  | 20.000 | ng/uI  | 0.00     |
| 79) Chrysene-d12              | 21.304 | 240  | 2008100  | 20.000 | ng/uI  | 0.00     |
| 88) Perylene-d12              | 23.686 | 264  | 1851318  | 20.000 | ng/uI  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.234  | 96   | 46951    | 6.091  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.658  | 84   | 655326   | 29.597 | ng/uI  | 0.00     |
| 7) Phenol-d5                  | 6.963  | 99   | 939901   | 33.765 | ng/uI  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 543181   | 32.687 | ng/uI  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.322  | 132  | 708055   | 34.832 | ng/uI  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.510  | 113  | 755932   | 34.336 | ng/uI  | 0.00     |
| 21) Nitrobenzene-d5           | 8.963  | 128  | 370720   | 40.198 | ng/uI  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.681  | 143  | 408195   | 42.643 | ng/uI  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.216 | 165  | 742488   | 35.791 | ng/uI  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.740 | 131  | 976743   | 31.622 | ng/uI  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.857 | 166  | 2187214  | 34.353 | ng/uI  | 0.00     |
| 49) Acenaphthylene-d8         | 14.134 | 160  | 2513966  | 34.552 | ng/uI  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.663 | 143  | 456549   | 37.888 | ng/uI  | 0.00     |
| 60) Fluorene-d10              | 15.440 | 176  | 1874580  | 34.389 | ng/uI  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.575 | 200  | 395486   | 38.167 | ng/uI  | 0.00     |
| 73) Anthracene-d10            | 17.304 | 188  | 2961858  | 34.539 | ng/uI  | 0.00     |
| 81) Pyrene-d10                | 19.545 | 212  | 3511163  | 34.826 | ng/uI  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.533 | 264  | 3238640  | 35.803 | ng/uI  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.270  | 88   | 99242    | 12.111 | ng/uL  | 98       |
| 5) Pyridine                   | 3.675  | 79   | 617020   | 28.356 | ng/uI  | 99       |
| 6) Benzaldehyde               | 6.940  | 77   | 459987   | 35.038 | ng/uI  | 100      |
| 8) Phenol                     | 6.993  | 94   | 936838   | 32.368 | ng/uI  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 7.222  | 93   | 732808   | 31.547 | ng/uI  | 99       |
| 12) 2-Chlorophenol            | 7.352  | 128  | 724628   | 32.746 | ng/uI  | 98       |
| 13) 2-Methylphenol            | 8.240  | 108  | 716179   | 32.176 | ng/uI  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.328  | 45   | 927890   | 29.786 | ng/uI  | 98       |
| 16) Acetophenone              | 8.622  | 105  | 1126899  | 32.956 | ng/uI  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.610  | 70   | 563615   | 34.238 | ng/uI  | 99       |
| 18) 4-Methylphenol            | 8.575  | 108  | 800668   | 32.912 | ng/uI  | 98       |
| 19) Hexachloroethane          | 8.863  | 117  | 282205   | 33.159 | ng/uI  | 97       |
| 22) Nitrobenzene              | 9.005  | 77   | 847899   | 35.238 | ng/uI  | 99       |
| 23) Isophorone                | 9.528  | 82   | 1657234  | 32.913 | ng/uI  | 100      |
| 25) 2-Nitrophenol             | 9.710  | 139  | 440067   | 38.341 | ng/uI  | 100      |
| 26) 2,4-Dimethylphenol        | 9.775  | 107  | 835062   | 33.134 | ng/uI  | 100      |
| 27) Bis(2-Chloroethoxy)met... | 10.010 | 93   | 1038569  | 32.232 | ng/uI  | 100      |
| 29) 2,4-Dichlorophenol        | 10.246 | 162  | 724141   | 33.626 | ng/uI  | 99       |
| 30) Naphthalene               | 10.640 | 128  | 2401126  | 31.857 | ng/uI  | 100      |
| 32) 4-Chloroaniline           | 10.763 | 127  | 973407   | 30.627 | ng/uI  | 99       |
| 33) Hexachlorobutadiene       | 10.916 | 225  | 422389   | 31.526 | ng/uI  | 98       |
| 34) Caprolactam               | 11.569 | 113  | 258634m  | 35.295 | ng/uI  |          |
| 35) 4-Chloro-3-methylphenol   | 11.899 | 107  | 807439   | 34.123 | ng/uI  | 99       |

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Manual IntegrationsAPPROVED

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| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.257 | 142  | 1655486  | 32.037 | ng/ul | 100      |
| 37) 1-Methylnaphthalene       | 12.475 | 142  | 1657054  | 32.110 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.628 | 216  | 855298   | 31.308 | ng/ul | 100      |
| 40) Hexachlorocyclopentadiene | 12.598 | 237  | 382805   | 30.288 | ng/ul | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.875 | 196  | 588017   | 33.673 | ng/ul | 99       |
| 42) 2,4,5-Trichlorophenol     | 12.946 | 196  | 646269   | 33.932 | ng/ul | 99       |
| 43) 1,1'-Biphenyl             | 13.275 | 154  | 2169563  | 31.724 | ng/ul | 100      |
| 44) 2-Chloronaphthalene       | 13.316 | 162  | 1711774  | 31.853 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.534 | 65   | 514324   | 41.293 | ng/ul | 99       |
| 47) Dimethylphthalate         | 13.904 | 163  | 2188940  | 32.632 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 14.034 | 165  | 455018   | 40.852 | ng/ul | 97       |
| 50) Acenaphthylene            | 14.163 | 152  | 2683023  | 32.882 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.363 | 138  | 491420   | 36.033 | ng/ul | 99       |
| 52) Acenaphthene              | 14.510 | 153  | 1836878  | 32.033 | ng/ul | 100      |
| 53) 2,4-Dinitrophenol         | 14.575 | 184  | 270605   | 35.976 | ng/ul | 96       |
| 55) 4-Nitrophenol             | 14.681 | 109  | 323733   | 35.507 | ng/ul | 96       |
| 56) Dibenzofuran              | 14.845 | 168  | 2535103  | 32.383 | ng/ul | 98       |
| 57) 2,4-Dinitrotoluene        | 14.828 | 165  | 666888   | 39.752 | ng/ul | 96       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.075 | 232  | 559070   | 34.584 | ng/ul | 97       |
| 59) Diethylphthalate          | 15.275 | 149  | 2178993  | 33.384 | ng/ul | 99       |
| 61) Fluorene                  | 15.498 | 166  | 2062090  | 33.152 | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.492 | 204  | 993570   | 32.565 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.534 | 138  | 521798m  | 42.349 | ng/ul |          |
| 66) 4,6-Dinitro-2-methylph... | 15.592 | 198  | 402999   | 36.503 | ng/ul | 97       |
| 67) N-Nitrosodiphenylamine    | 15.710 | 169  | 1804540  | 32.193 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.392 | 248  | 658694   | 32.950 | ng/ul | 99       |
| 69) Hexachlorobenzene         | 16.498 | 284  | 787116   | 32.829 | ng/ul | 99       |
| 70) Atrazine                  | 16.675 | 200  | 562079   | 28.946 | ng/ul | 100      |
| 71) Pentachlorophenol         | 16.857 | 266  | 486027   | 33.007 | ng/ul | 99       |
| 72) Phenanthrene              | 17.245 | 178  | 3450657  | 32.819 | ng/ul | 100      |
| 74) Anthracene                | 17.339 | 178  | 3430805  | 32.846 | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.234 | 216  | 904530   | 31.771 | ng/uL | 100      |
| 76) Pentachlorobenzene        | 14.763 | 250  | 872490   | 29.021 | ng/uL | 99       |
| 77) Carbazole                 | 17.616 | 167  | 3284997  | 35.311 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 18.163 | 149  | 3803338  | 35.191 | ng/ul | 100      |
| 80) Fluoranthene              | 19.222 | 202  | 4149635  | 33.336 | ng/ul | 99       |
| 82) Pyrene                    | 19.575 | 202  | 4228469  | 32.796 | ng/ul | 99       |
| 83) Butylbenzylphthalate      | 20.445 | 149  | 1832405  | 38.120 | ng/ul | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.222 | 252  | 1436221  | 32.925 | ng/ul | 100      |
| 85) Benzo(a)anthracene        | 21.286 | 228  | 4227254  | 32.997 | ng/ul | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.204 | 149  | 2617080  | 35.606 | ng/ul | 99       |
| 87) Chrysene                  | 21.339 | 228  | 3974751  | 33.188 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.122 | 149  | 4479811  | 40.569 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 22.963 | 252  | 4086660  | 34.860 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.004 | 252  | 4093215  | 36.637 | ng/ul | 100      |
| 93) Benzo(a)pyrene            | 23.586 | 252  | 3491443  | 34.105 | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.168 | 276  | 3936470  | 30.858 | ng/ul | 100      |
| 95) Dibenzo(a,h)anthracene    | 26.180 | 278  | 3400118  | 29.767 | ng/ul | 100      |
| 96) Benzo(g,h,i)perylene      | 26.933 | 276  | 3223891  | 28.589 | ng/ul | 99       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SLCS596

**Manual IntegrationsAPPROVED**

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## Manual IntegrationsAPPROVED

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