

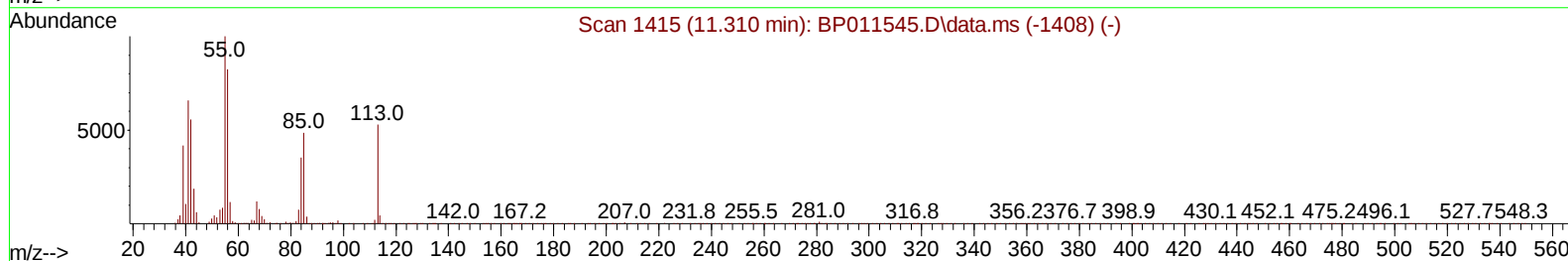
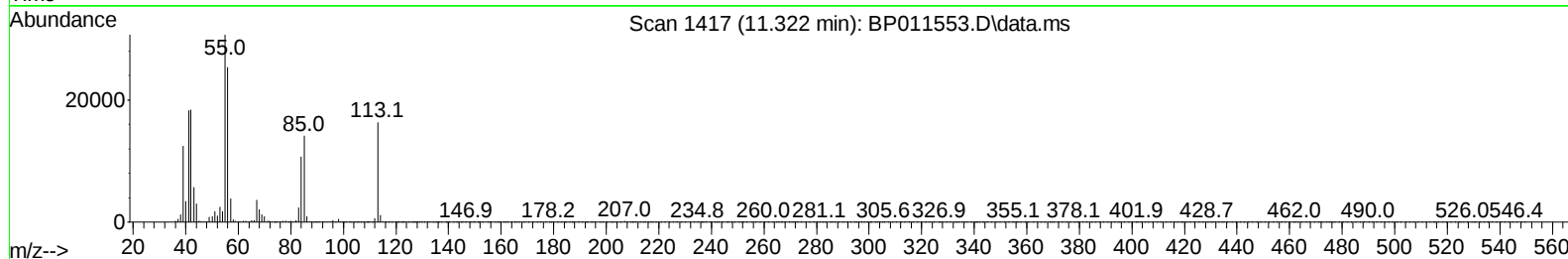
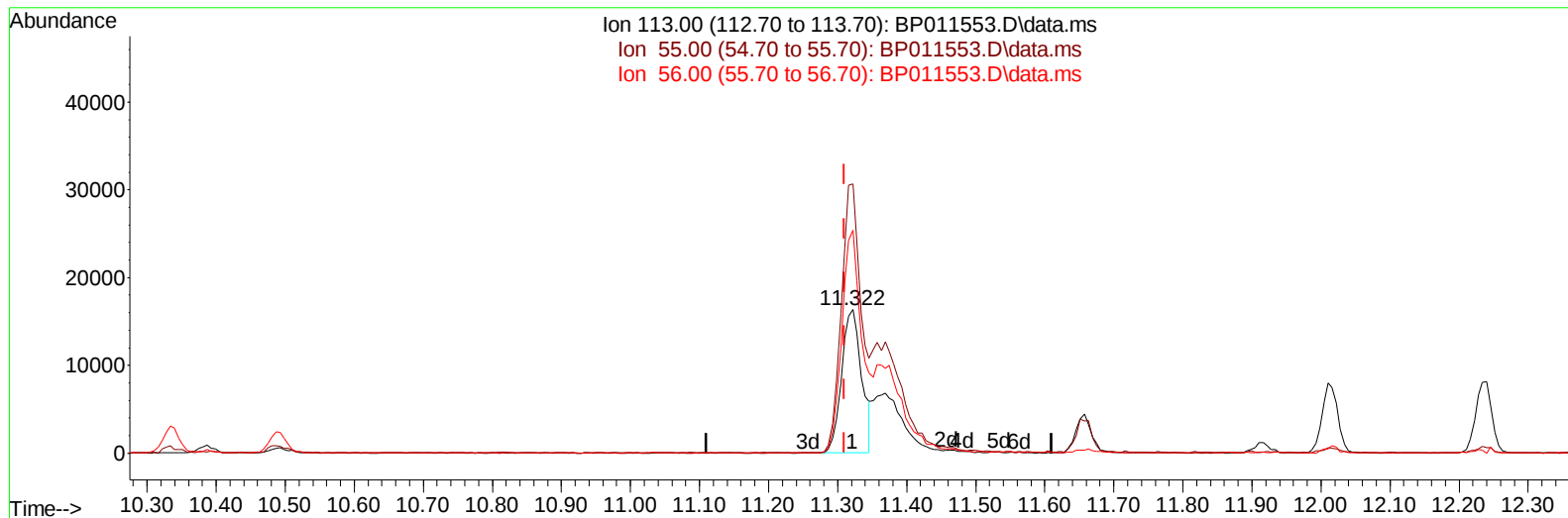
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
Data File : BP011553.D  
Acq On : 24 Aug 2022 22:30  
Operator : CG/JU  
Sample : PB147243BS  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS243

Manual IntegrationsAPPROVED

Quant Time: Aug 25 02:20:06 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 25 02:14:54 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/25/2022  
Supervised By :mohammad ahmed 08/27/2022



TIC: BP011553.D\data.ms

(34) Caprolactam

11.322min (+ 0.012) 21.24 ng/ul

response 33076

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 191.80 | 187.43 |
| 56.00  | 149.50 | 155.28 |
| 0.00   | 0.00   | 0.00   |

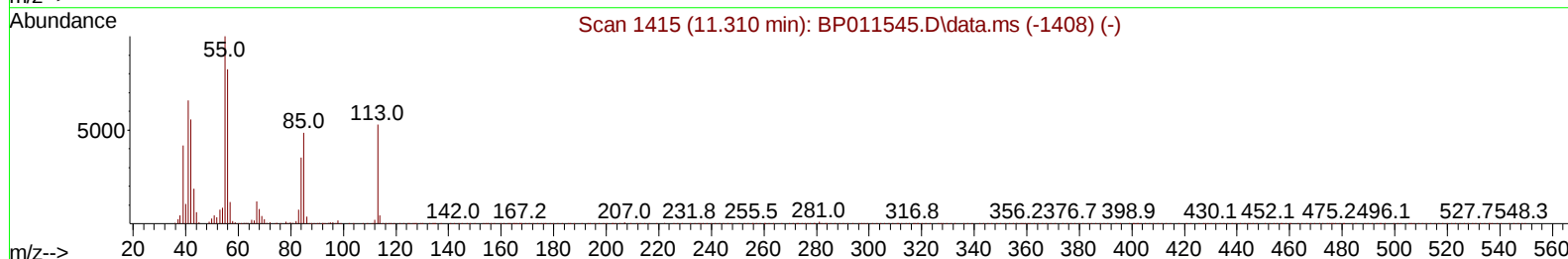
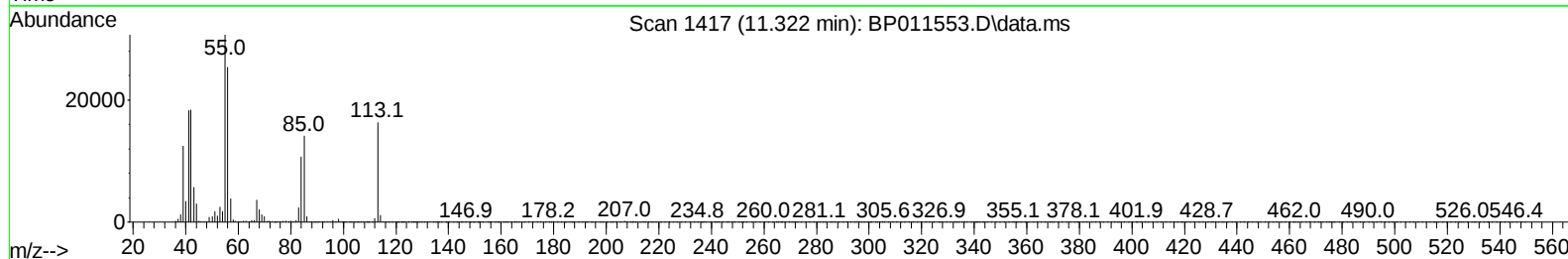
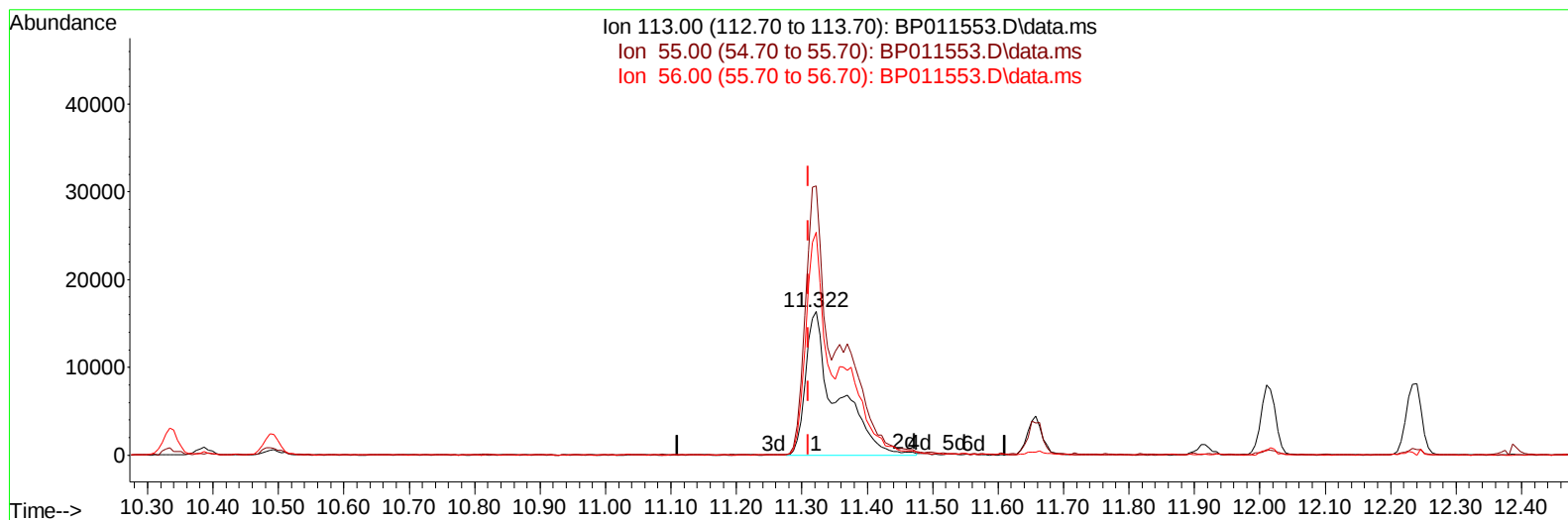
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
Data File : BP011553.D  
Acq On : 24 Aug 2022 22:30  
Operator : CG/JU  
Sample : PB147243BS  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS243

Manual IntegrationsAPPROVED

Quant Time: Aug 25 02:20:06 2022  
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Quant Title : SVOA CALIBRATION  
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Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/25/2022  
Supervised By :mohammad ahmed 08/27/2022



TIC: BP011553.D\data.ms

**(34) Caprolactam**

**11.322min (+ 0.012) 34.72 ng/ul m**

**response 54058**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 191.80 | 187.43 |
| 56.00  | 149.50 | 155.28 |
| 0.00   | 0.00   | 0.00   |

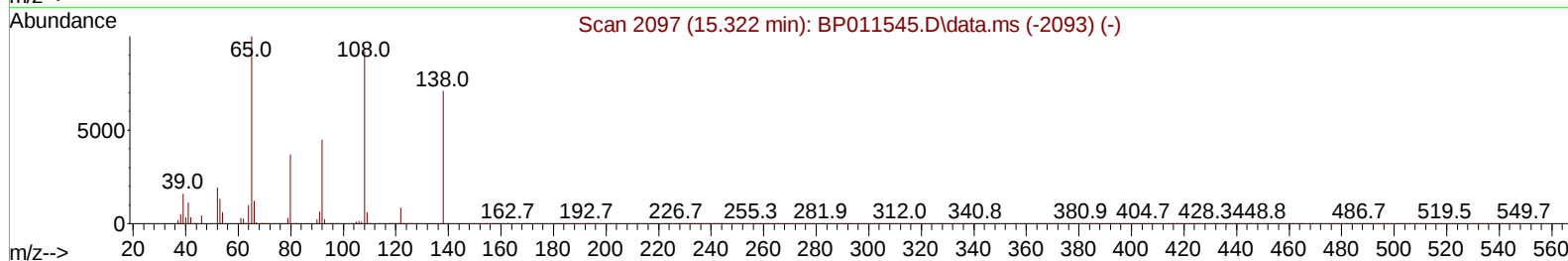
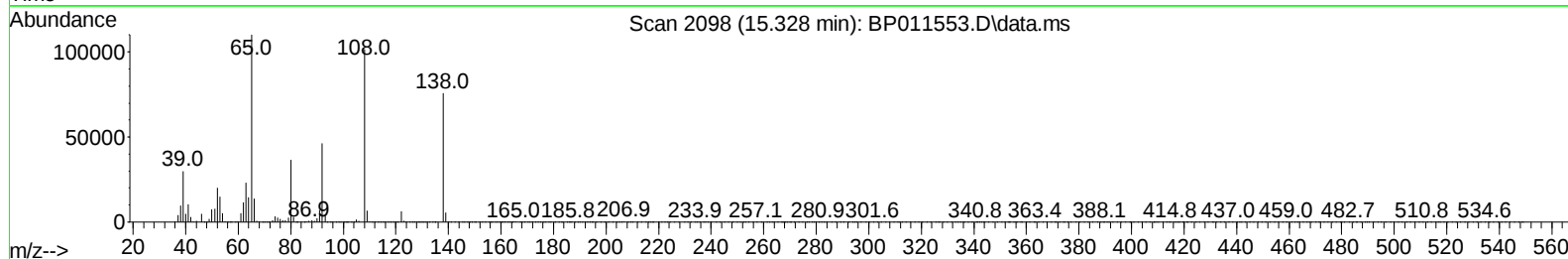
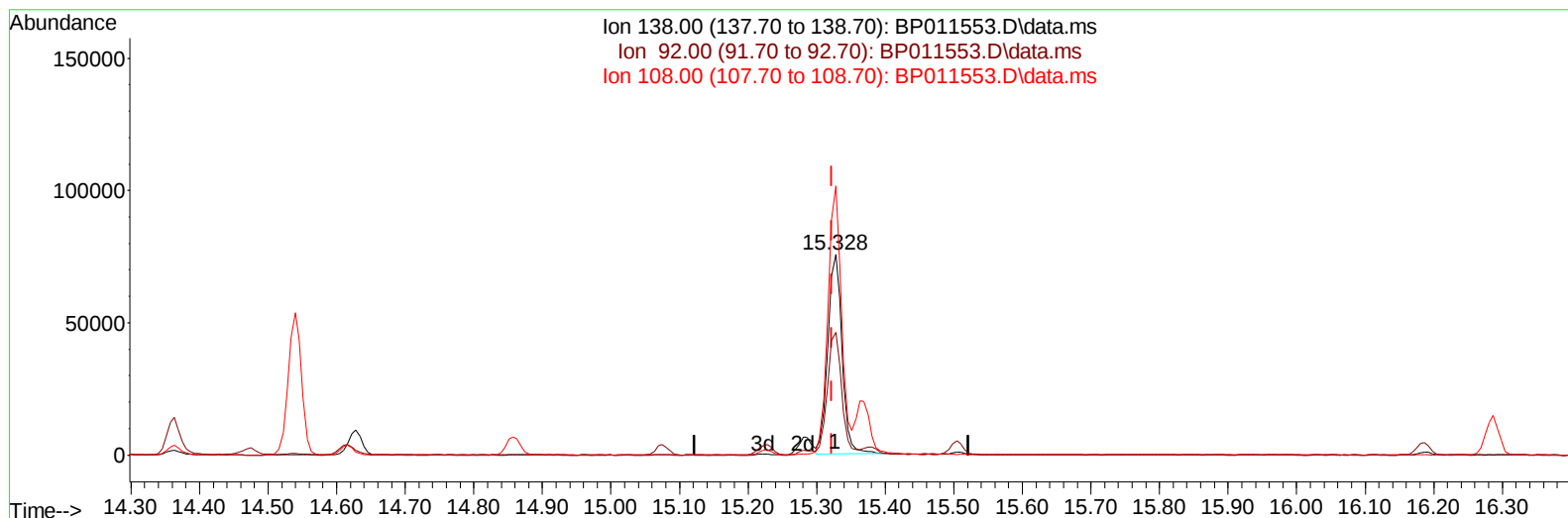
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
Data File : BP011553.D  
Acq On : 24 Aug 2022 22:30  
Operator : CG/JU  
Sample : PB147243BS  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS243

Manual IntegrationsAPPROVED

Quant Time: Aug 25 02:20:06 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 25 02:14:54 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/25/2022  
Supervised By :mohammad ahmed 08/27/2022



TIC: BP011553.D\data.ms

**(63) 4-Nitroaniline**

**15.328min (+ 0.006) 40.72 ng/ul**

**response 109076**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 61.90  | 61.02  |
| 108.00 | 117.10 | 133.96 |
| 0.00   | 0.00   | 0.00   |

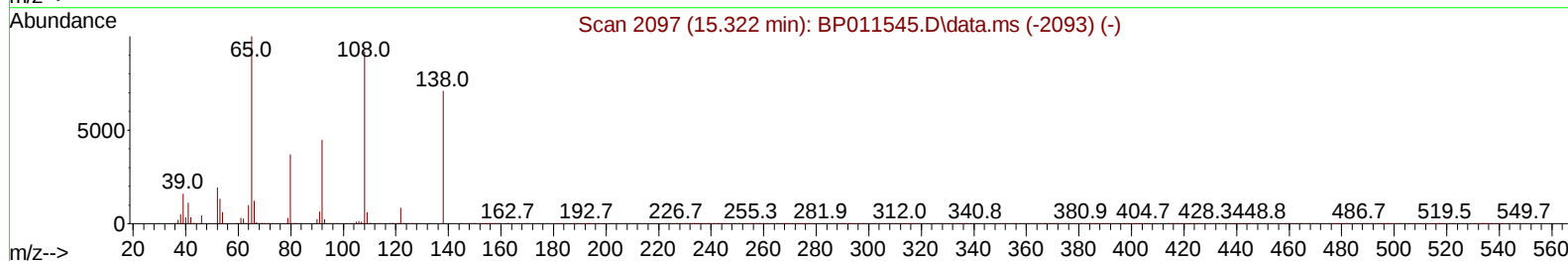
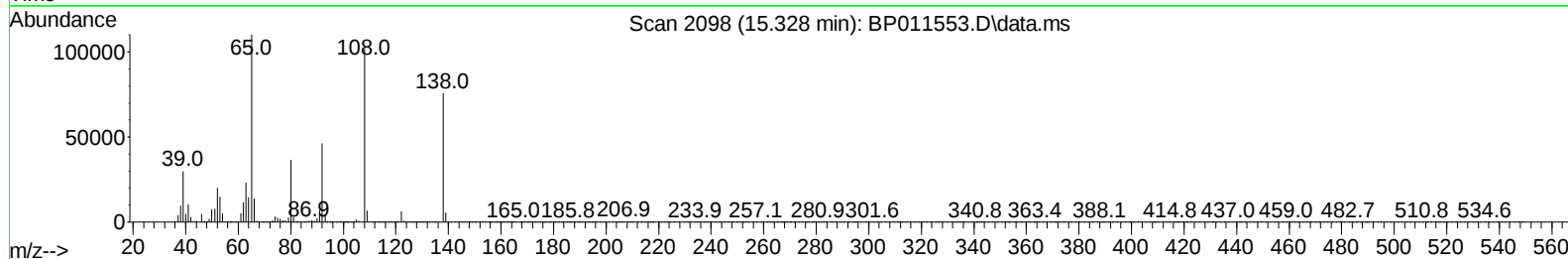
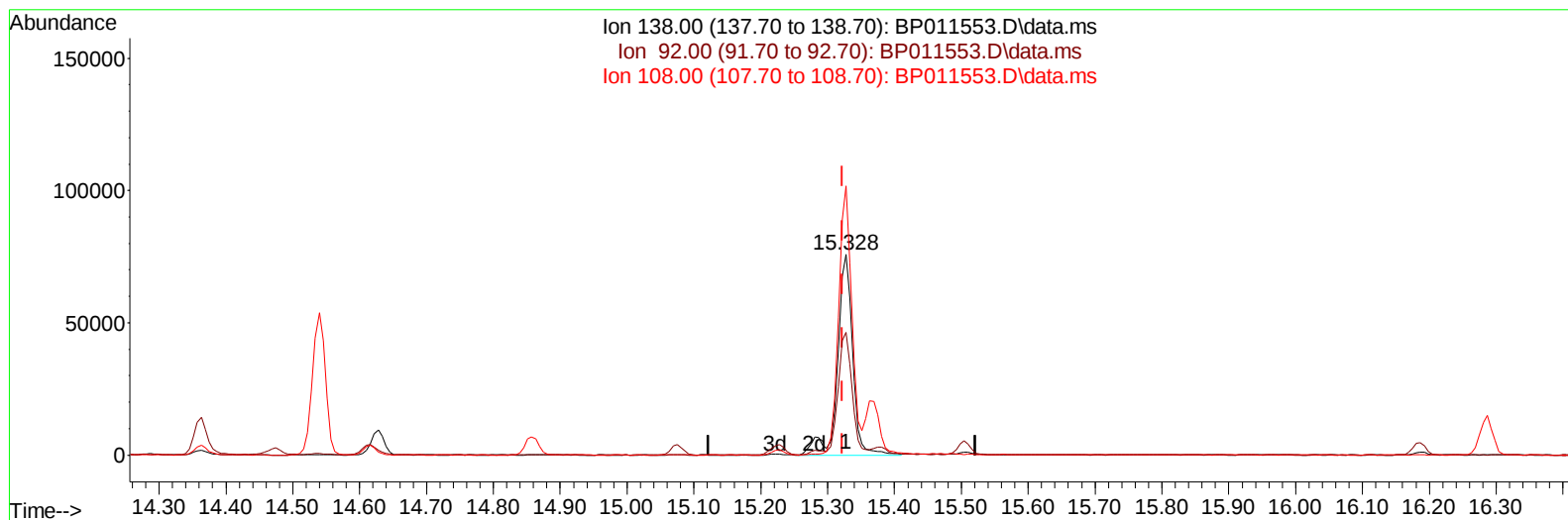
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
 Data File : BP011553.D  
 Acq On : 24 Aug 2022 22:30  
 Operator : CG/JU  
 Sample : PB147243BS  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS243

Manual IntegrationsAPPROVED

Quant Time: Aug 25 02:20:06 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 25 02:14:54 2022  
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Reviewed By :Jagrut Upadhyay 08/25/2022  
 Supervised By :mohammad ahmed 08/27/2022



TIC: BP011553.D\data.ms

**(63) 4-Nitroaniline**

**15.328min (+ 0.006) 45.41 ng/ul m**

**response 121643**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 61.90  | 61.02  |
| 108.00 | 117.10 | 133.96 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
 Data File : BP011553.D  
 Acq On : 24 Aug 2022 22:30  
 Operator : CG/JU  
 Sample : PB147243BS  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS243

Manual IntegrationsAPPROVED

Quant Time: Aug 25 02:20:06 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 25 02:14:54 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/25/2022  
 Supervised By :mohammad ahmed 08/27/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.546  | 152  | 69712    | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.334 | 136  | 310260   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.222 | 164  | 193727   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 16.992 | 188  | 394306   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.122 | 240  | 399894   | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 23.410 | 264  | 403853   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.064  | 96   | 13834    | 6.158  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.470  | 84   | 170263   | 29.603 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.734  | 99   | 211529   | 33.019 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.899  | 67   | 135447   | 32.544 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.081  | 132  | 167712   | 35.594 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.275  | 113  | 171020   | 33.353 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.716  | 128  | 83116    | 37.211 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.428  | 143  | 88244    | 38.819 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.963  | 165  | 157148   | 37.451 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.493 | 131  | 225880   | 32.338 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.651 | 166  | 496572   | 36.241 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.910 | 160  | 559838   | 36.762 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.457 | 143  | 99535    | 37.703 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.228 | 176  | 415100   | 36.010 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.369 | 200  | 79943    | 37.521 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.092 | 188  | 667720   | 38.236 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.357 | 212  | 751550   | 37.146 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.263 | 264  | 770500   | 39.221 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.099  | 88   | 26684    | 11.392 | ng/uL  | 99       |
| 5) Pyridine                   | 3.493  | 79   | 164689   | 28.686 | ng/ul  | 97       |
| 6) Benzaldehyde               | 6.705  | 77   | 89453    | 30.427 | ng/ul  | 97       |
| 8) Phenol                     | 6.763  | 94   | 204367   | 31.043 | ng/ul  | 96       |
| 10) Bis(2-Chloroethyl)ether   | 6.993  | 93   | 158627   | 30.889 | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.116  | 128  | 169180   | 33.368 | ng/ul  | 97       |
| 13) 2-Methylphenol            | 8.005  | 108  | 153308   | 31.611 | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.093  | 45   | 241005   | 31.679 | ng/ul  | 99       |
| 16) Acetophenone              | 8.381  | 105  | 267835   | 31.400 | ng/ul  | 97       |
| 17) N-Nitroso-di-n-propyla... | 8.375  | 70   | 139595   | 32.901 | ng/ul  | 99       |
| 18) 4-Methylphenol            | 8.340  | 108  | 169683   | 31.602 | ng/ul  | 98       |
| 19) Hexachloroethane          | 8.610  | 117  | 78473    | 34.121 | ng/ul  | 97       |
| 22) Nitrobenzene              | 8.757  | 77   | 214992   | 34.147 | ng/ul  | 98       |
| 23) Isophorone                | 9.287  | 82   | 380170   | 34.399 | ng/ul  | 100      |
| 25) 2-Nitrophenol             | 9.463  | 139  | 93603    | 36.089 | ng/ul  | 93       |
| 26) 2,4-Dimethylphenol        | 9.534  | 107  | 211225   | 34.299 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 9.775  | 93   | 221916   | 33.219 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 9.993  | 162  | 151795   | 35.347 | ng/ul  | 99       |
| 30) Naphthalene               | 10.387 | 128  | 537230   | 32.794 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.516 | 127  | 166053   | 22.945 | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 10.663 | 225  | 95375    | 34.887 | ng/ul  | 93       |
| 34) Caprolactam               | 11.322 | 113  | 54058m   | 34.717 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 11.657 | 107  | 194629   | 35.847 | ng/ul  | 97       |

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 Sample : PB147243BS  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLC5243

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/25/2022  
 Supervised By :mohammad ahmed 08/27/2022

Quant Time: Aug 25 02:20:06 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 25 02:14:54 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.016 | 142  | 359617   | 33.741 | ng/ul | 99       |
| 37) 1-Methylnaphthalene       | 12.234 | 142  | 363193   | 33.598 | ng/ul | 96       |
| 39) 1,2,4,5-Tetrachloroben... | 12.387 | 216  | 167480   | 34.061 | ng/ul | 98       |
| 40) Hexachlorocyclopentadiene | 12.357 | 237  | 89036    | 29.187 | ng/ul | 97       |
| 41) 2,4,6-Trichlorophenol     | 12.640 | 196  | 110800   | 35.011 | ng/ul | 100      |
| 42) 2,4,5-Trichlorophenol     | 12.716 | 196  | 122021   | 34.637 | ng/ul | 99       |
| 43) 1,1'-Biphenyl             | 13.045 | 154  | 480076   | 33.258 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.087 | 162  | 360134   | 33.108 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.310 | 65   | 140463   | 37.835 | ng/ul | 98       |
| 47) Dimethylphthalate         | 13.698 | 163  | 488937   | 34.525 | ng/ul | 98       |
| 48) 2,6-Dinitrotoluene        | 13.822 | 165  | 96181    | 37.776 | ng/ul | 98       |
| 50) Acenaphthylene            | 13.940 | 152  | 616774   | 34.091 | ng/ul | 99       |
| 51) 3-Nitroaniline            | 14.151 | 138  | 105234   | 37.946 | ng/ul | 98       |
| 52) Acenaphthene              | 14.287 | 153  | 409547   | 33.797 | ng/ul | 99       |
| 53) 2,4-Dinitrophenol         | 14.363 | 184  | 55458    | 35.587 | ng/ul | 94       |
| 55) 4-Nitrophenol             | 14.475 | 109  | 119326   | 37.936 | ng/ul | 91       |
| 56) Dibenzofuran              | 14.628 | 168  | 568898   | 34.044 | ng/ul | 98       |
| 57) 2,4-Dinitrotoluene        | 14.616 | 165  | 148650   | 39.000 | ng/ul | 90       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.857 | 232  | 95966    | 34.697 | ng/ul | 97       |
| 59) Diethylphthalate          | 15.075 | 149  | 563267   | 36.314 | ng/ul | 98       |
| 61) Fluorene                  | 15.281 | 166  | 473240   | 34.617 | ng/ul | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.287 | 204  | 216349   | 35.032 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.328 | 138  | 121643m  | 45.409 | ng/ul |          |
| 66) 4,6-Dinitro-2-methylph... | 15.381 | 198  | 81115    | 36.804 | ng/ul | 93       |
| 67) N-Nitrosodiphenylamine    | 15.504 | 169  | 414564   | 35.836 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.186 | 248  | 126187   | 35.850 | ng/ul | 96       |
| 69) Hexachlorobenzene         | 16.286 | 284  | 148070   | 36.501 | ng/ul | 98       |
| 70) Atrazine                  | 16.475 | 200  | 46213    | 10.704 | ng/ul | 98       |
| 71) Pentachlorophenol         | 16.645 | 266  | 83141    | 33.054 | ng/ul | 98       |
| 72) Phenanthrene              | 17.039 | 178  | 759492   | 35.586 | ng/ul | 99       |
| 74) Anthracene                | 17.128 | 178  | 782935   | 36.362 | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.004 | 216  | 169615   | 33.463 | ng/uL | 99       |
| 76) Pentachlorobenzene        | 14.540 | 250  | 166003   | 34.515 | ng/uL | 98       |
| 77) Carbazole                 | 17.410 | 167  | 727360   | 35.914 | ng/ul | 99       |
| 78) Di-n-butylphthalate       | 17.980 | 149  | 994094   | 39.859 | ng/ul | 99       |
| 80) Fluoranthene              | 19.028 | 202  | 893792   | 35.922 | ng/ul | 98       |
| 82) Pyrene                    | 19.386 | 202  | 919891   | 35.186 | ng/ul | 97       |
| 83) Butylbenzylphthalate      | 20.286 | 149  | 471951   | 40.242 | ng/ul | 94       |
| 84) 3,3'-Dichlorobenzidine    | 21.051 | 252  | 280864   | 37.100 | ng/ul | 99       |
| 85) Benzo(a)anthracene        | 21.110 | 228  | 890983   | 36.570 | ng/ul | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.051 | 149  | 720837   | 41.534 | ng/ul | 98       |
| 87) Chrysene                  | 21.163 | 228  | 883839   | 36.929 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 21.945 | 149  | 1222189  | 37.648 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 22.716 | 252  | 925630   | 36.846 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 22.763 | 252  | 879935   | 36.705 | ng/ul | 98       |
| 93) Benzo(a)pyrene            | 23.316 | 252  | 896168   | 36.539 | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 25.768 | 276  | 1023991  | 37.432 | ng/ul | 100      |
| 95) Dibenzo(a,h)anthracene    | 25.786 | 278  | 882381   | 37.460 | ng/ul | 97       |
| 96) Benzo(g,h,i)perylene      | 26.492 | 276  | 820211   | 35.248 | ng/ul | 99       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SLCS243

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 08/25/2022  
Supervised By :mohammad ahmed 08/27/2022

Manual IntegrationsAPPROVED

Quant Time: Aug 25 02:20:06 2022  
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