

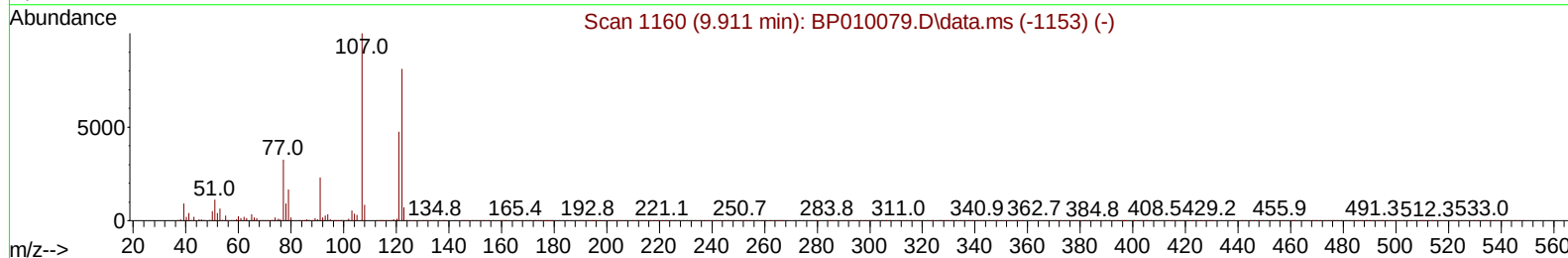
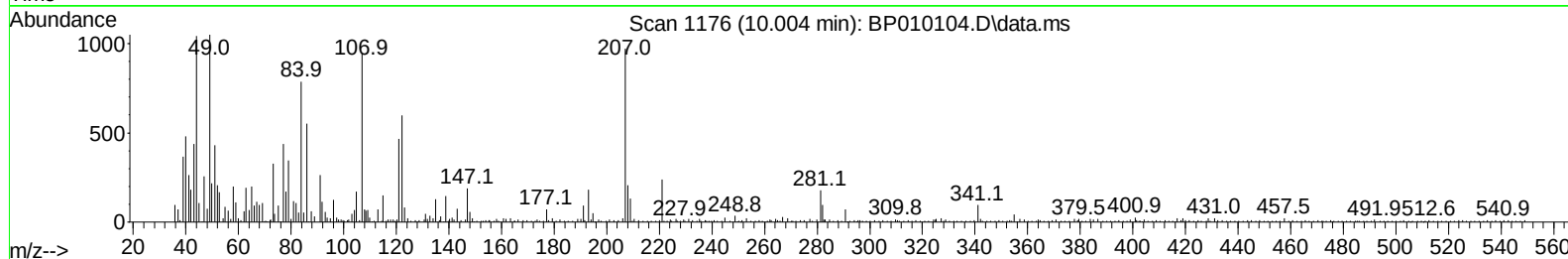
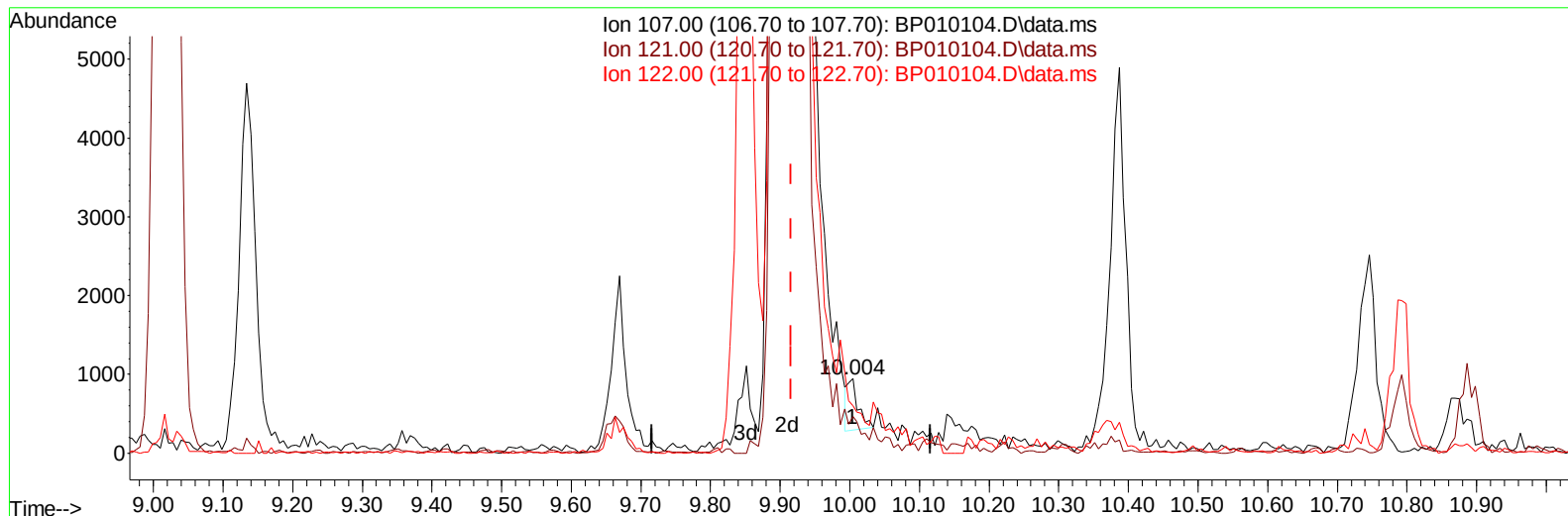
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010104.D\data.ms

(26) 2,4-Dimethylphenol

10.004min (+ 0.088) 0.04 ng/ul

response 688

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 107.00 | 100.00 | 100.00 |
| 121.00 | 42.60  | 49.26  |
| 122.00 | 57.90  | 63.35  |
| 0.00   | 0.00   | 0.00   |

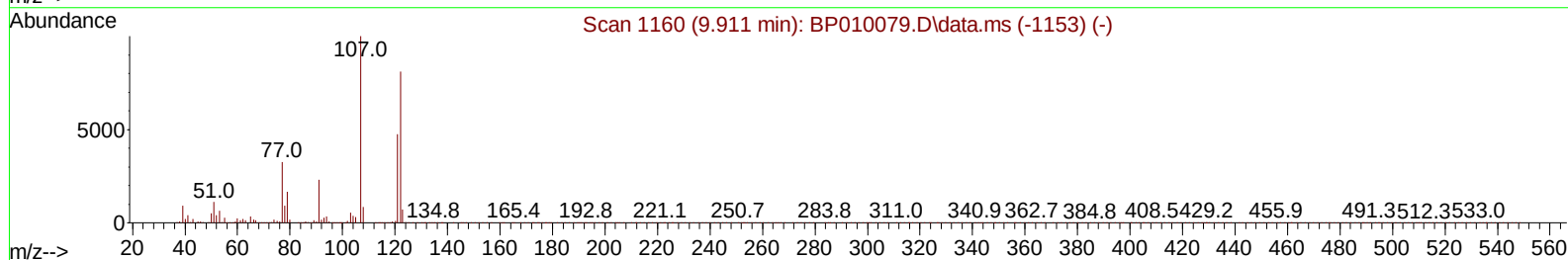
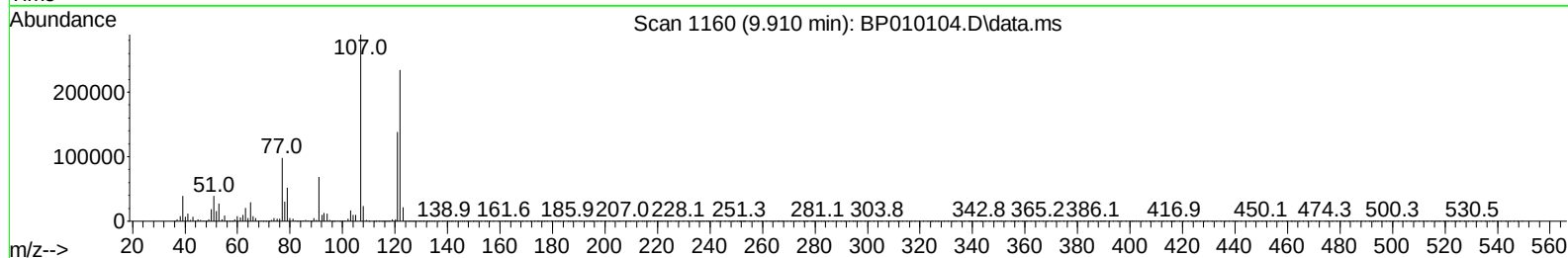
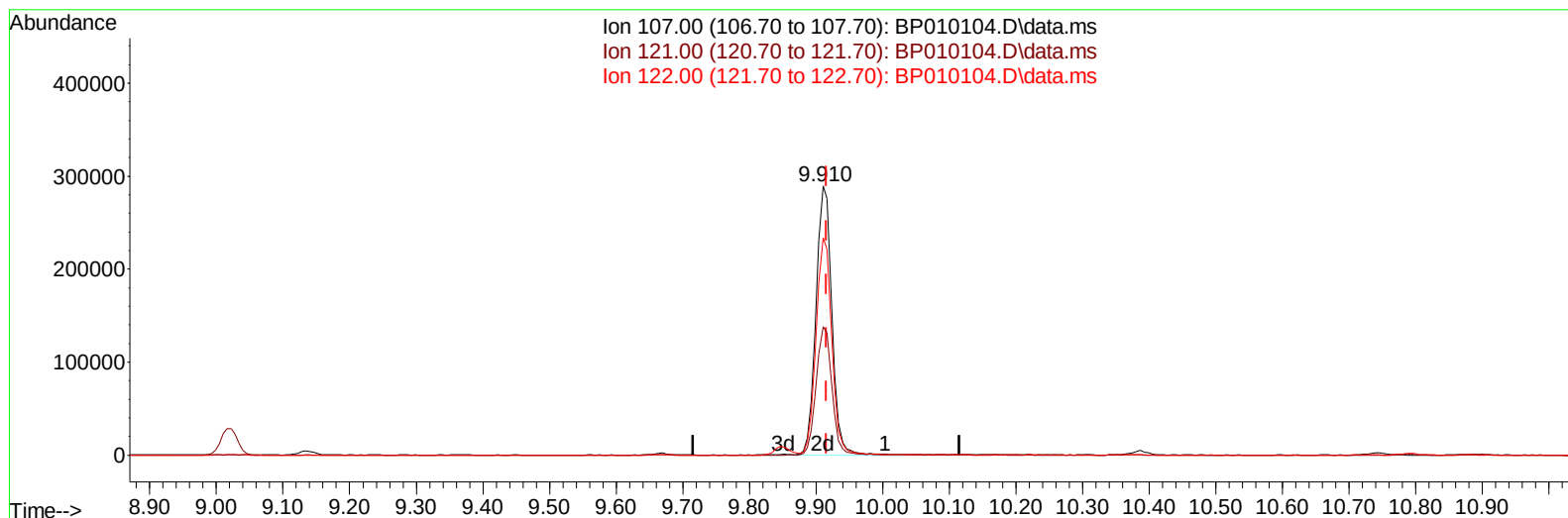
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010104.D\data.ms

(26) 2,4-Dimethylphenol

9.910min (-0.006) 30.35 ng/ul m

response 477688

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 107.00 | 100.00 | 100.00 |
| 121.00 | 42.60  | 47.78  |
| 122.00 | 57.90  | 80.86# |
| 0.00   | 0.00   | 0.00   |

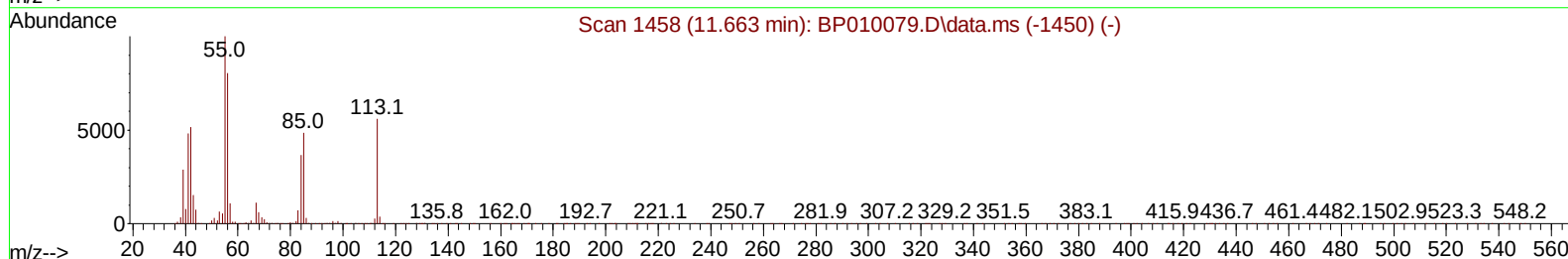
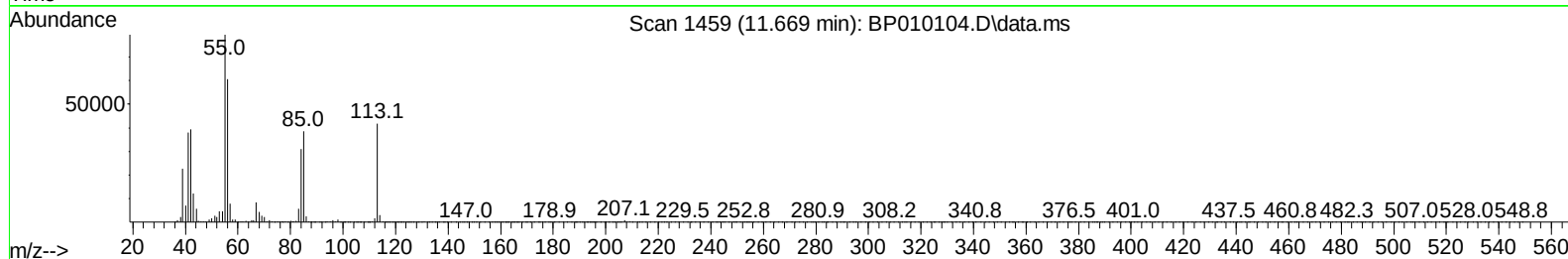
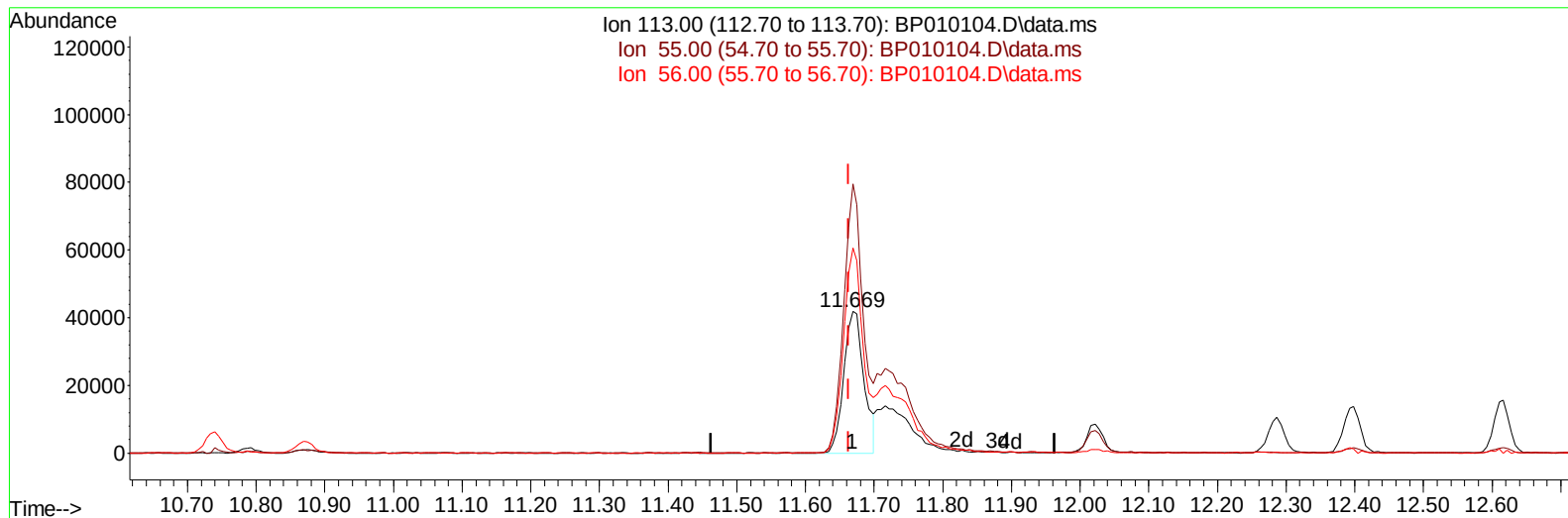
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010104.D\data.ms

**(34) Caprolactam**

**11.669min (+ 0.006) 18.58 ng/ul**

**response 85852**

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00  |
| 55.00  | 72.70  | 189.92# |
| 56.00  | 85.80  | 144.77# |
| 0.00   | 0.00   | 0.00    |

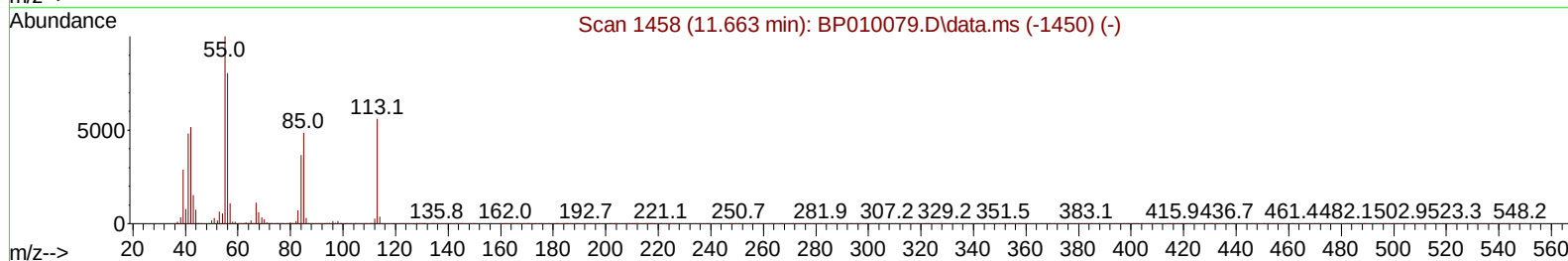
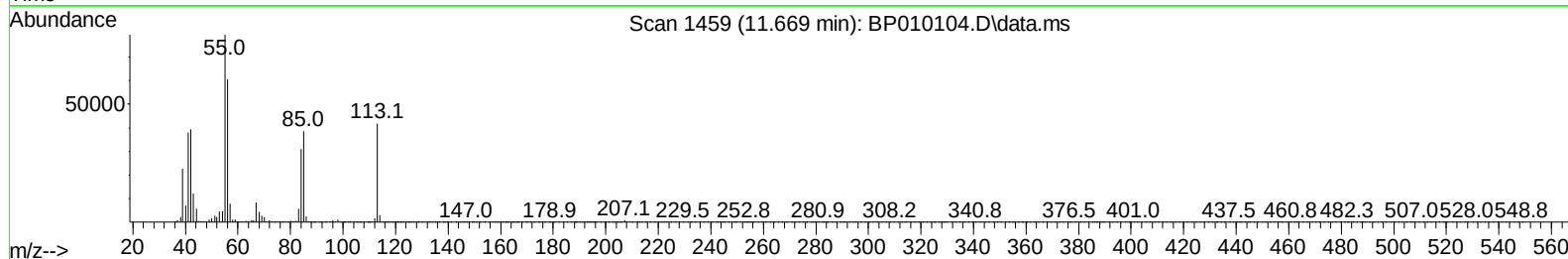
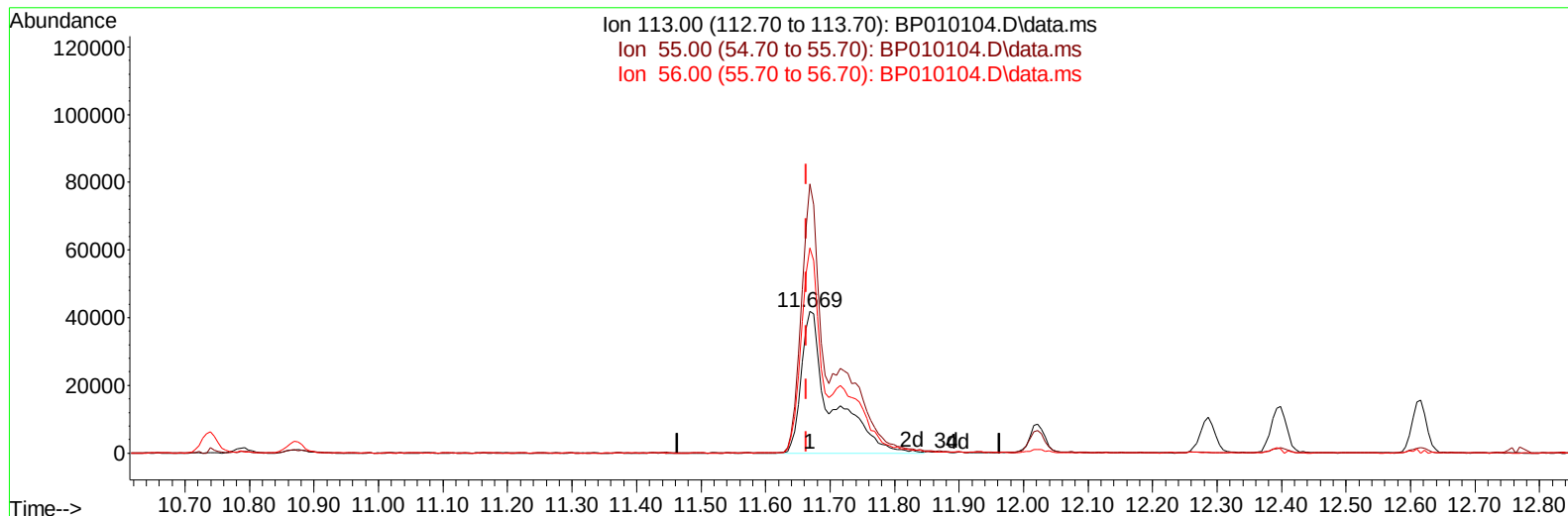
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
Data File : BP010104.D  
Acq On : 26 Apr 2022 03:11  
Operator : CG/JU  
Sample : PB144328BS  
Misc :  
ALS Vial : 26 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Apr 21 14:49:38 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
Supervised By :Yogesh Patel 04/28/2022



TIC: BP010104.D\data.ms

(34) Caprolactam

11.669min (+ 0.006) 29.30 ng/ul m

response 135400

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00  |
| 55.00  | 72.70  | 189.92# |
| 56.00  | 85.80  | 144.77# |
| 0.00   | 0.00   | 0.00    |

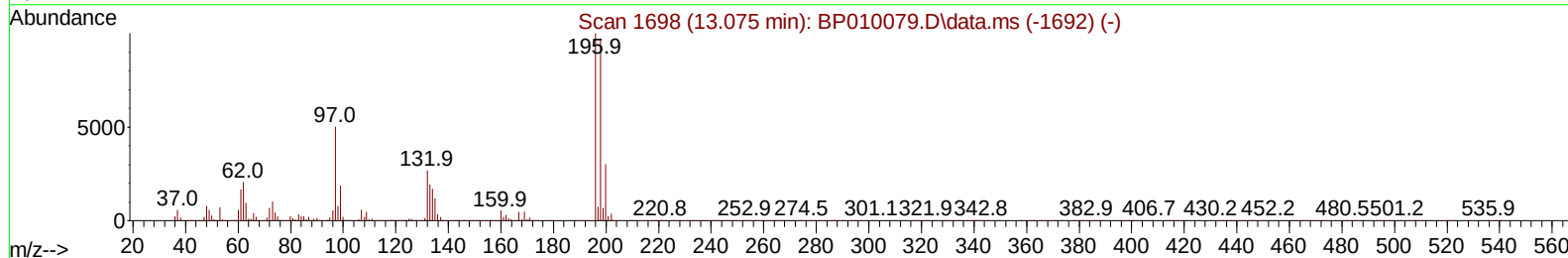
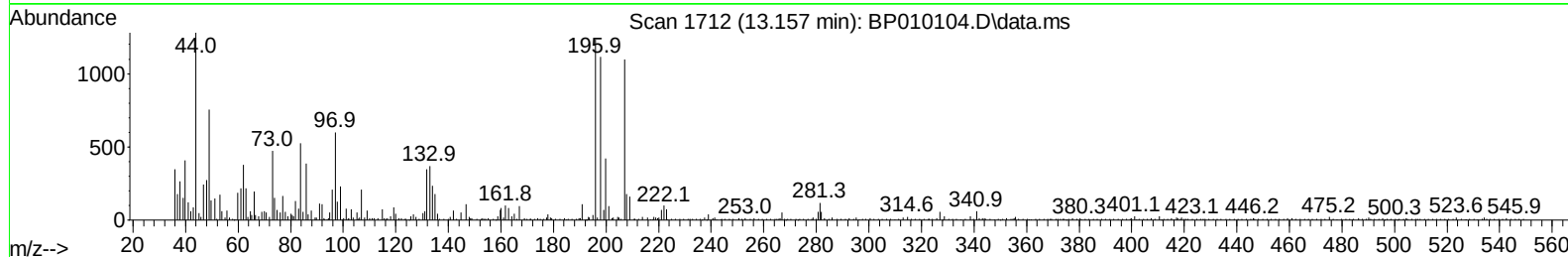
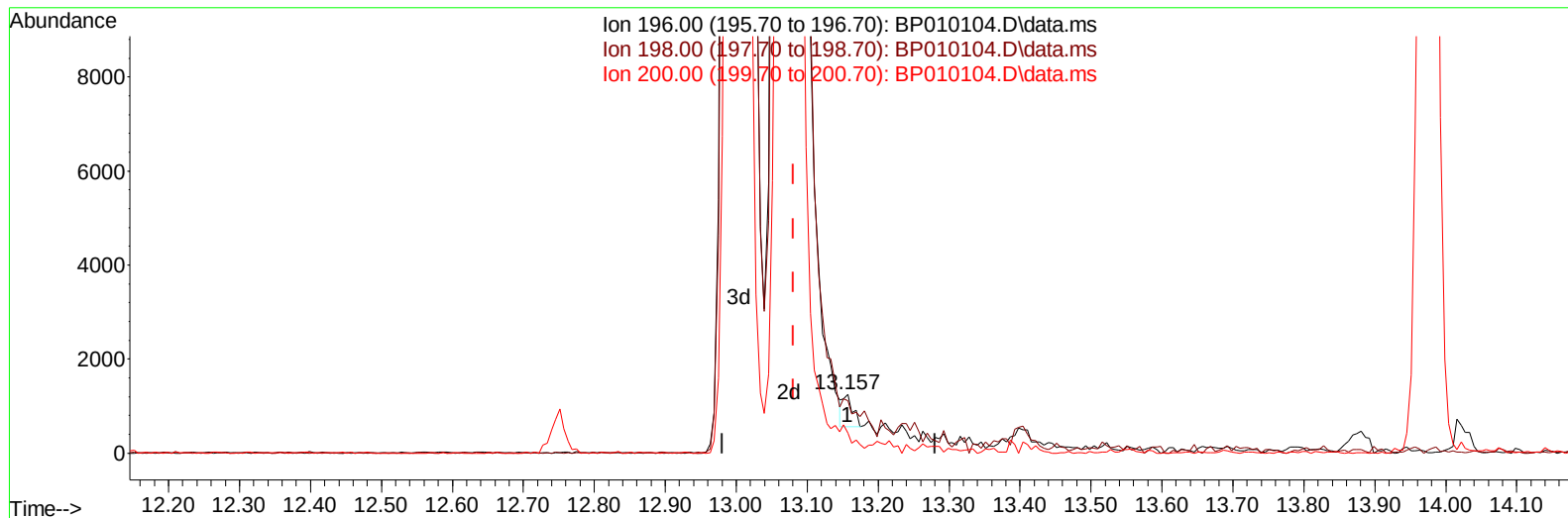
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010104.D\data.ms

(42) 2,4,5-Trichlorophenol

13.157min (+ 0.076) 0.06 ng/ul

response 672

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 196.00 | 100.00 | 100.00 |
| 198.00 | 85.80  | 89.35  |
| 200.00 | 41.00  | 33.87  |
| 0.00   | 0.00   | 0.00   |

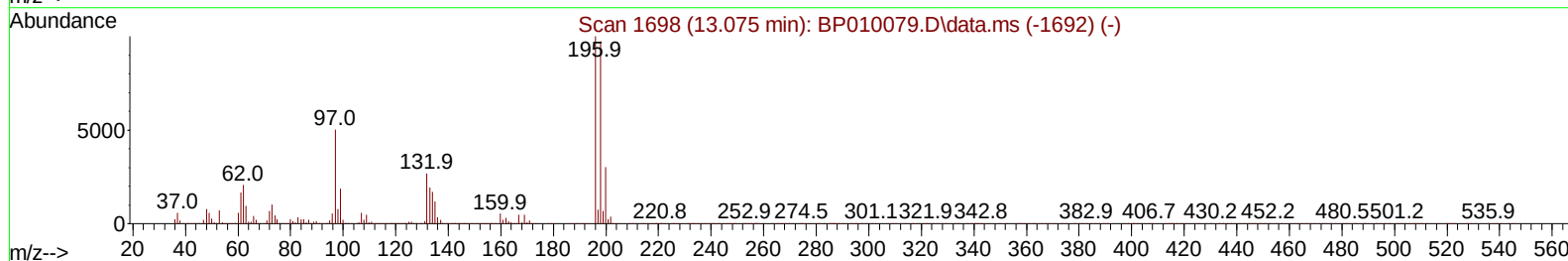
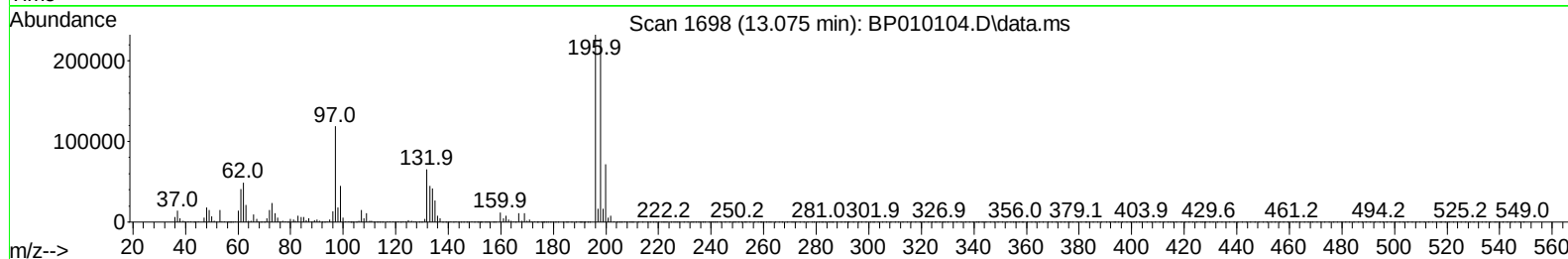
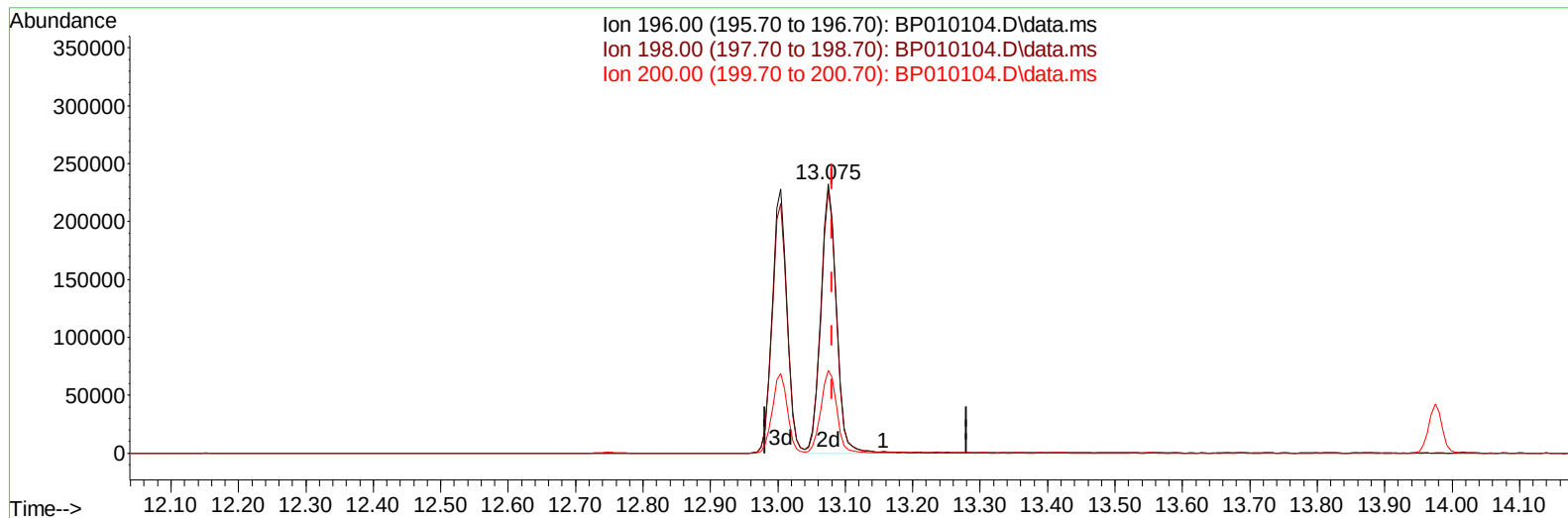
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
Data File : BP010104.D  
Acq On : 26 Apr 2022 03:11  
Operator : CG/JU  
Sample : PB144328BS  
Misc :  
ALS Vial : 26 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Apr 21 14:49:38 2022  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
Supervised By :Yogesh Patel 04/28/2022



TIC: BP010104.D\data.ms

(42) 2,4,5-Trichlorophenol

13.075min (-0.006) 32.73 ng/ul m

response 375986

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 196.00 | 100.00 | 100.00 |
| 198.00 | 85.80  | 97.99  |
| 200.00 | 41.00  | 30.76# |
| 0.00   | 0.00   | 0.00   |

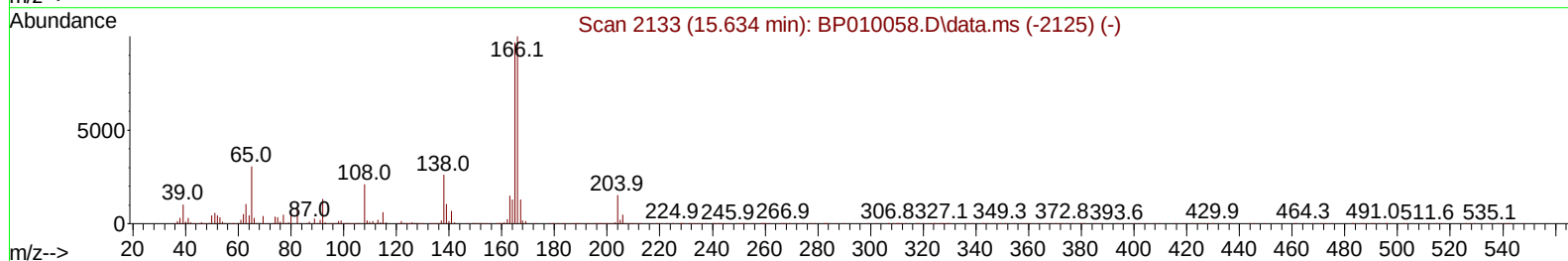
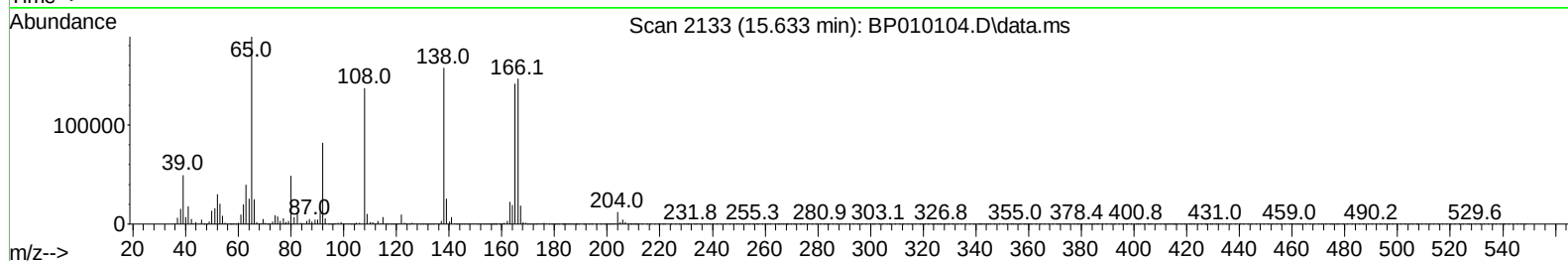
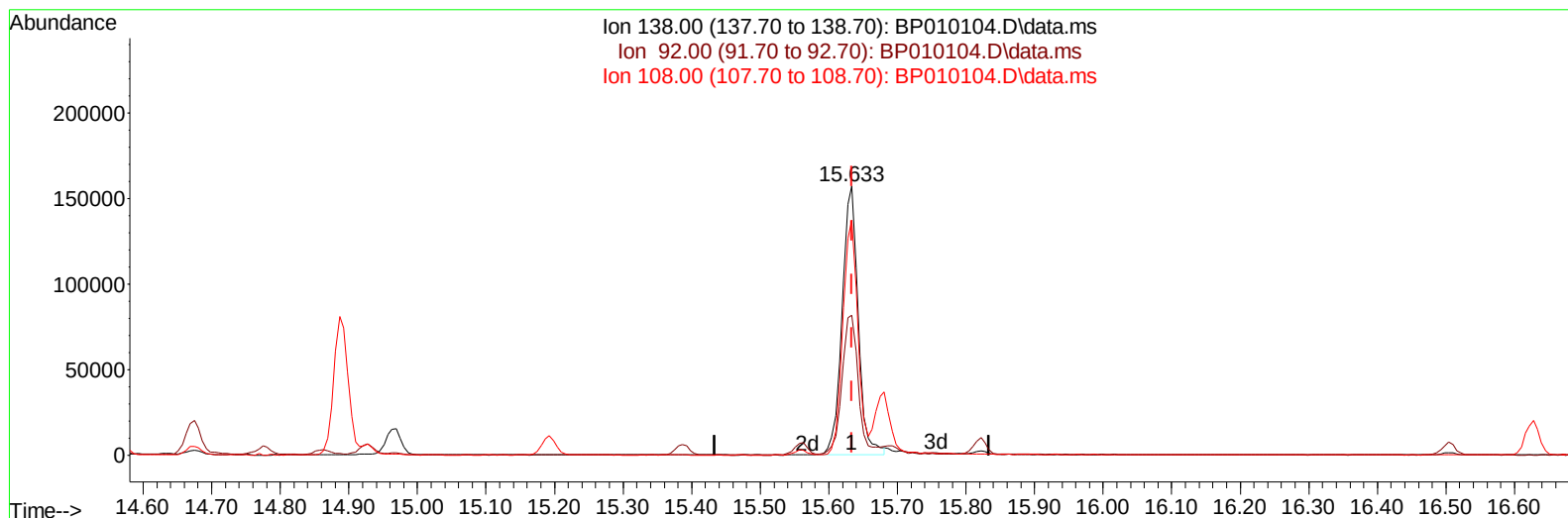
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010104.D\data.ms

**(63) 4-Nitroaniline**

**15.633min (-0.000) 35.01 ng/ul**

**response 252816**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 52.00  | 51.95  |
| 108.00 | 122.00 | 87.31# |
| 0.00   | 0.00   | 0.00   |

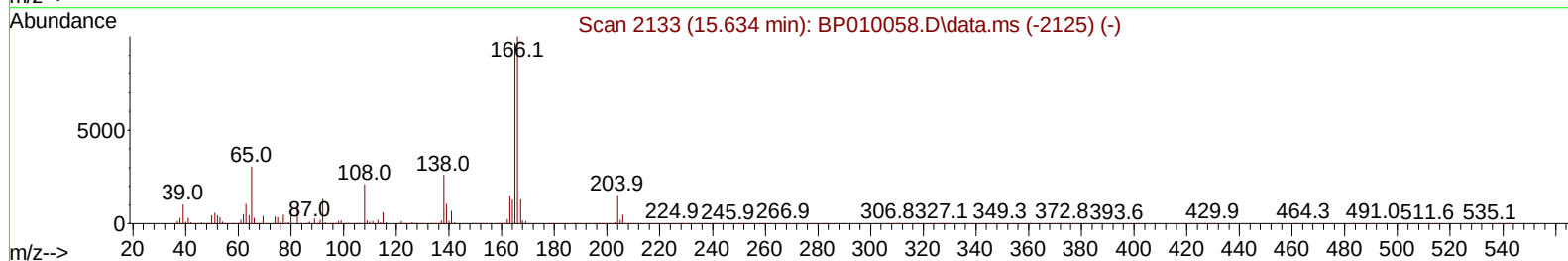
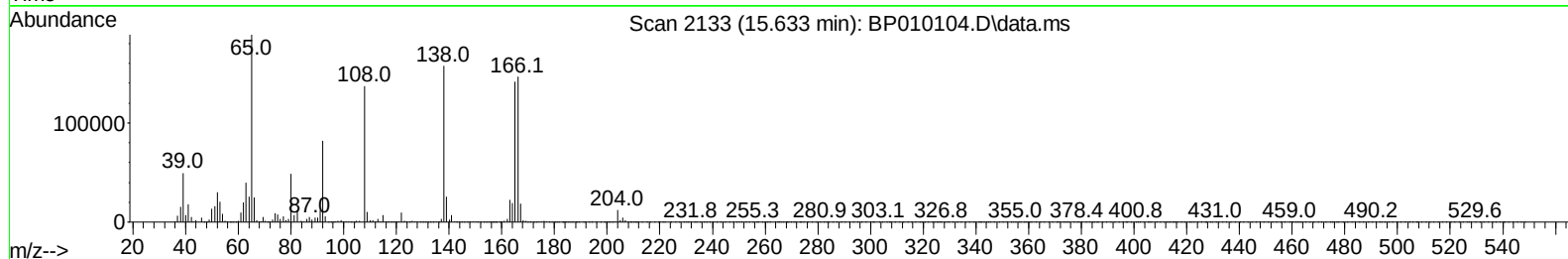
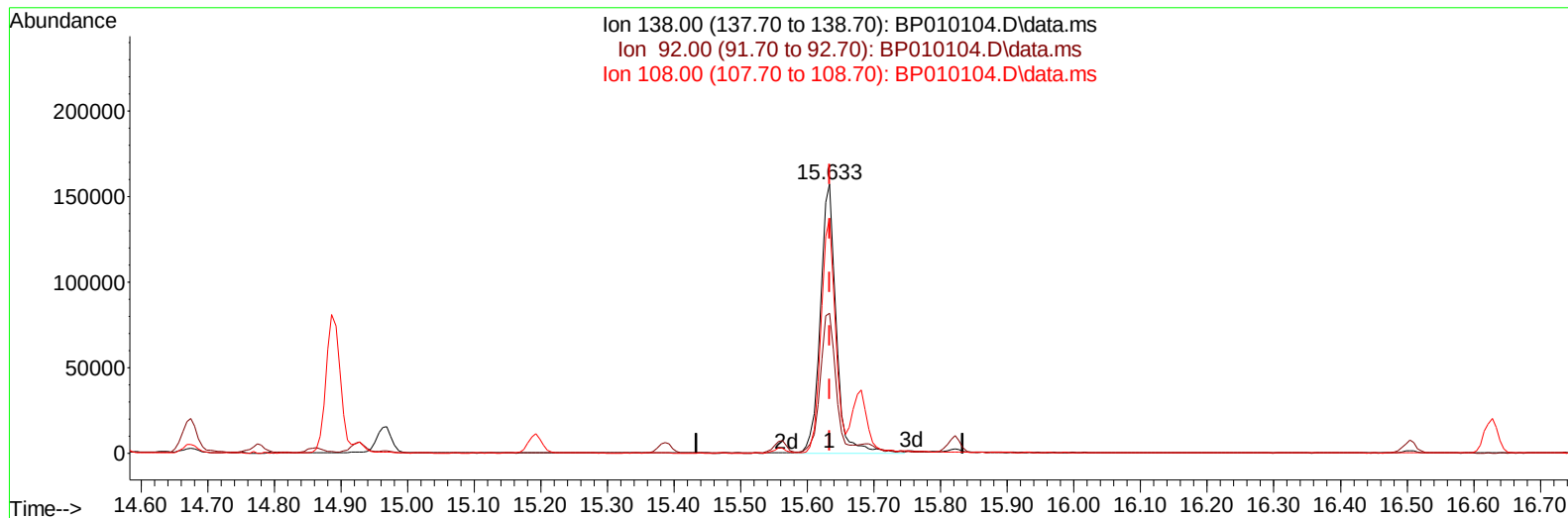
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010104.D\data.ms

**(63) 4-Nitroaniline**

**15.633min (-0.000) 36.26 ng/ul m**

**response 261869**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 138.00 | 100.00 | 100.00 |
| 92.00  | 52.00  | 51.95  |
| 108.00 | 122.00 | 87.31# |
| 0.00   | 0.00   | 0.00   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.934  | 152  | 185345   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.740 | 136  | 790204   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.569 | 164  | 528104   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.316 | 188  | 1123531  | 20.000 | ng/ul  | # 0.00   |
| 79) Chrysene-d12              | 21.404 | 240  | 976579   | 20.000 | ng/ul  | # 0.00   |
| 88) Perylene-d12              | 23.851 | 264  | 735355   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.369  | 96   | 24234    | 5.702  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 319889   | 26.753 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.099  | 99   | 475835   | 28.225 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.263  | 67   | 309188   | 30.564 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.463  | 132  | 383470   | 30.713 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.646  | 113  | 409507   | 29.317 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.093  | 128  | 198589   | 31.598 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.816  | 143  | 223308   | 32.534 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.357 | 165  | 406779   | 30.442 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.869 | 131  | 488654   | 25.505 | ng/ul  | -0.01    |
| 46) Dimethylphthalate-d6      | 13.975 | 166  | 1282681  | 31.224 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.257 | 160  | 1481391  | 29.795 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.763 | 143  | 209120   | 27.651 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.557 | 176  | 1091889  | 30.899 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.681 | 200  | 213976   | 30.119 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.422 | 188  | 1673458  | 30.787 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.651 | 212  | 1936298  | 34.864 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.692 | 264  | 1269709  | 32.935 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.399  | 88   | 48270    | 10.929 | ng/ul# | 16       |
| 5) Pyridine                   | 3.799  | 79   | 340717   | 27.412 | ng/ul# | 45       |
| 6) Benzaldehyde               | 7.069  | 77   | 335223   | 37.208 | ng/ul  | 95       |
| 8) Phenol                     | 7.122  | 94   | 501039   | 29.578 | ng/ul# | 93       |
| 10) Bis(2-Chloroethyl)ether   | 7.352  | 93   | 394306   | 30.213 | ng/ul# | 79       |
| 12) 2-Chlorophenol            | 7.499  | 128  | 384127   | 30.501 | ng/ul  | 94       |
| 13) 2-Methylphenol            | 8.375  | 108  | 381095   | 29.334 | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.463  | 45   | 582165   | 30.062 | ng/ul# | 85       |
| 16) Acetophenone              | 8.757  | 105  | 625362   | 28.696 | ng/ul# | 81       |
| 17) N-Nitroso-di-n-propyla... | 8.746  | 70   | 332202   | 29.095 | ng/ul# | 73       |
| 18) 4-Methylphenol            | 8.710  | 108  | 416531   | 29.120 | ng/ul  | 97       |
| 19) Hexachloroethane          | 9.022  | 117  | 163569   | 30.909 | ng/ul  | 83       |
| 22) Nitrobenzene              | 9.134  | 77   | 482036   | 31.304 | ng/ul# | 84       |
| 23) Isophorone                | 9.669  | 82   | 934099   | 30.713 | ng/ul# | 96       |
| 25) 2-Nitrophenol             | 9.851  | 139  | 232246   | 31.947 | ng/ul# | 83       |
| 26) 2,4-Dimethylphenol        | 9.910  | 107  | 477688m  | 30.351 | ng/ul  |          |
| 27) Bis(2-Chloroethoxy)met... | 10.146 | 93   | 549672   | 30.835 | ng/ul# | 97       |
| 29) 2,4-Dichlorophenol        | 10.387 | 162  | 402561   | 31.340 | ng/ul# | 82       |
| 30) Naphthalene               | 10.793 | 128  | 1285596  | 31.084 | ng/ul  | 97       |
| 32) 4-Chloroaniline           | 10.898 | 127  | 474837   | 25.150 | ng/ul  | 97       |
| 33) Hexachlorobutadiene       | 11.081 | 225  | 285673   | 31.380 | ng/ul  | 95       |
| 34) Caprolactam               | 11.669 | 113  | 135400m  | 29.305 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.022 | 107  | 466182   | 30.043 | ng/ul# | 69       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.398 | 142  | 904816   | 30.165 | ng/ul  | 90       |
| 37) 1-Methylnaphthalene       | 12.616 | 142  | 924173   | 30.344 | ng/ul# | 95       |
| 39) 1,2,4,5-Tetrachloroben... | 12.769 | 216  | 530778   | 32.531 | ng/ul  | 95       |
| 40) Hexachlorocyclopentadiene | 12.751 | 237  | 304550   | 33.122 | ng/ul  | 95       |
| 41) 2,4,6-Trichlorophenol     | 13.004 | 196  | 347202   | 32.889 | ng/ul# | 85       |
| 42) 2,4,5-Trichlorophenol     | 13.075 | 196  | 375986m  | 32.734 | ng/ul  |          |
| 43) 1,1'-Biphenyl             | 13.404 | 154  | 1256246  | 32.252 | ng/ul  | 93       |
| 44) 2-Chloronaphthalene       | 13.445 | 162  | 968068   | 32.116 | ng/ul  | 94       |
| 45) 2-Nitroaniline            | 13.645 | 65   | 315821   | 32.312 | ng/ul# | 82       |
| 47) Dimethylphthalate         | 14.022 | 163  | 1232259  | 30.864 | ng/ul# | 92       |
| 48) 2,6-Dinitrotoluene        | 14.139 | 165  | 262166   | 31.997 | ng/ul  | 94       |
| 50) Acenaphthylene            | 14.287 | 152  | 1553503  | 31.544 | ng/ul  | 95       |
| 51) 3-Nitroaniline            | 14.469 | 138  | 230343   | 31.034 | ng/ul# | 81       |
| 52) Acenaphthene              | 14.634 | 153  | 992162   | 30.983 | ng/ul  | 97       |
| 53) 2,4-Dinitrophenol         | 14.675 | 184  | 123819   | 27.044 | ng/ul# | 78       |
| 55) 4-Nitrophenol             | 14.775 | 109  | 182237   | 27.735 | ng/ul# | 47       |
| 56) Dibenzofuran              | 14.963 | 168  | 1462500  | 30.842 | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene        | 14.928 | 165  | 378391   | 31.613 | ng/ul# | 85       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.192 | 232  | 315188   | 30.871 | ng/ul# | 86       |
| 59) Diethylphthalate          | 15.386 | 149  | 1269689  | 31.266 | ng/ul  | 93       |
| 61) Fluorene                  | 15.616 | 166  | 1188175  | 30.518 | ng/ul  | 91       |
| 62) 4-Chlorophenyl-phenyle... | 15.610 | 204  | 624506   | 30.479 | ng/ul  | 99       |
| 63) 4-Nitroaniline            | 15.633 | 138  | 261869m  | 36.259 | ng/ul  |          |
| 66) 4,6-Dinitro-2-methylph... | 15.692 | 198  | 208203   | 30.376 | ng/ul# | 87       |
| 67) N-Nitrosodiphenylamine    | 15.822 | 169  | 1036933  | 32.166 | ng/ul  | 94       |
| 68) 4-Bromophenyl-phenylether | 16.504 | 248  | 392017   | 32.711 | ng/ul  | 95       |
| 69) Hexachlorobenzene         | 16.628 | 284  | 442771   | 31.830 | ng/ul# | 91       |
| 70) Atrazine                  | 16.780 | 200  | 413500   | 31.550 | ng/ul# | 88       |
| 71) Pentachlorophenol         | 16.969 | 266  | 267735   | 30.223 | ng/ul# | 84       |
| 72) Phenanthrene              | 17.363 | 178  | 1896834  | 31.062 | ng/ul  | 98       |
| 74) Anthracene                | 17.451 | 178  | 1895950  | 30.824 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.369 | 216  | 541258   | 32.999 | ng/ul# | 86       |
| 76) Pentachlorobenzene        | 14.892 | 250  | 511584   | 31.806 | ng/ul  | 91       |
| 77) Carbazole                 | 17.722 | 167  | 1646723  | 30.254 | ng/ul# | 96       |
| 78) Di-n-butylphthalate       | 18.269 | 149  | 2172769  | 32.447 | ng/ul# | 93       |
| 80) Fluoranthene              | 19.322 | 202  | 2245118  | 35.840 | ng/ul  | 97       |
| 82) Pyrene                    | 19.680 | 202  | 2265556  | 35.103 | ng/ul# | 93       |
| 83) Butylbenzylphthalate      | 20.545 | 149  | 946863   | 35.492 | ng/ul# | 81       |
| 84) 3,3'-Dichlorobenzidine    | 21.316 | 252  | 609865   | 34.707 | ng/ul# | 93       |
| 85) Benzo(a)anthracene        | 21.386 | 228  | 1970337  | 31.792 | ng/ul  | 95       |
| 86) Bis(2-ethylhexyl)phtha... | 21.310 | 149  | 1396166  | 35.310 | ng/ul# | 82       |
| 87) Chrysene                  | 21.439 | 228  | 1957996  | 32.085 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.251 | 149  | 2241100  | 39.203 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.104 | 252  | 1665358  | 33.687 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.157 | 252  | 1615747  | 34.504 | ng/ul# | 97       |
| 93) Benzo(a)pyrene            | 23.745 | 252  | 1509310  | 33.015 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.415 | 276  | 1451711  | 31.922 | ng/ul# | 92       |
| 95) Dibenzo(a,h)anthracene    | 26.427 | 278  | 1239806  | 31.860 | ng/ul# | 96       |
| 96) Benzo(g,h,i)perylene      | 27.198 | 276  | 1245810  | 32.466 | ng/ul  | 93       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SLCS328

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 04/26/2022  
Supervised By :Yogesh Patel 04/28/2022

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042522\  
 Data File : BP010104.D  
 Acq On : 26 Apr 2022 03:11  
 Operator : CG/JU  
 Sample : PB144328BS  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS328

Manual IntegrationsAPPROVED

Quant Time: Apr 26 04:18:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 14:49:38 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/28/2022

