

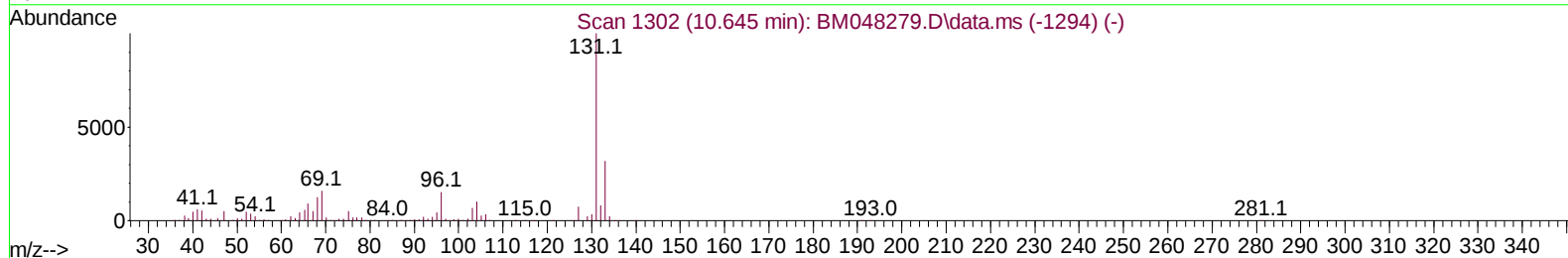
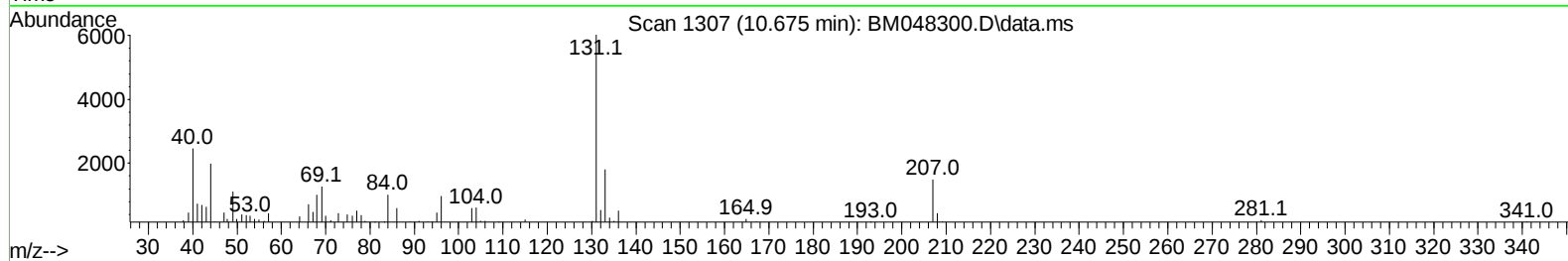
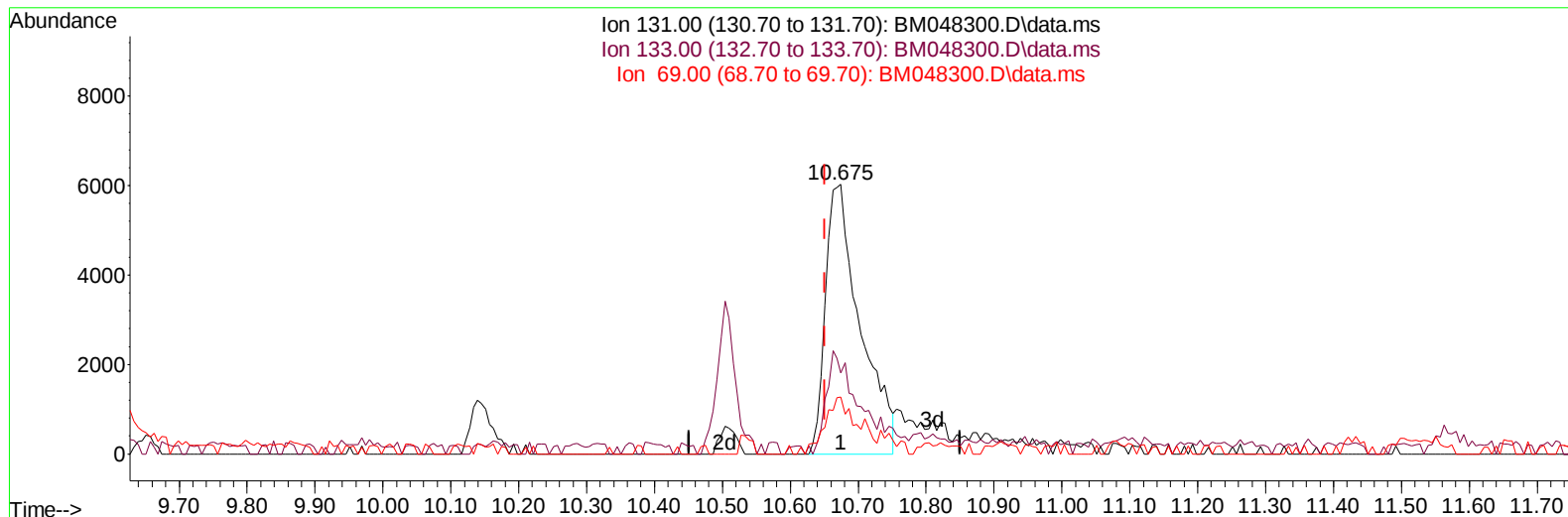
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM102824\  
Data File : BM048300.D  
Acq On : 29 Oct 2024 23:18  
Operator : RC/JU  
Sample : P4492-11  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
A4EQ6

Manual IntegrationsAPPROVED

Quant Time: Oct 29 23:51:42 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM102624.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun Oct 27 23:39:58 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/30/2024  
Supervised By :mohammad ahmed 11/01/2024



TIC: BM048300.D\data.ms

**(31) 4-Chloroaniline-d4 (S)**

**10.675min (+ 0.024) 1.30 ng/ul**

**response 21374**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 131.00 | 100.00 | 100.00 |
| 133.00 | 31.80  | 30.09  |
| 69.00  | 16.30  | 21.05# |
| 0.00   | 0.00   | 0.00   |

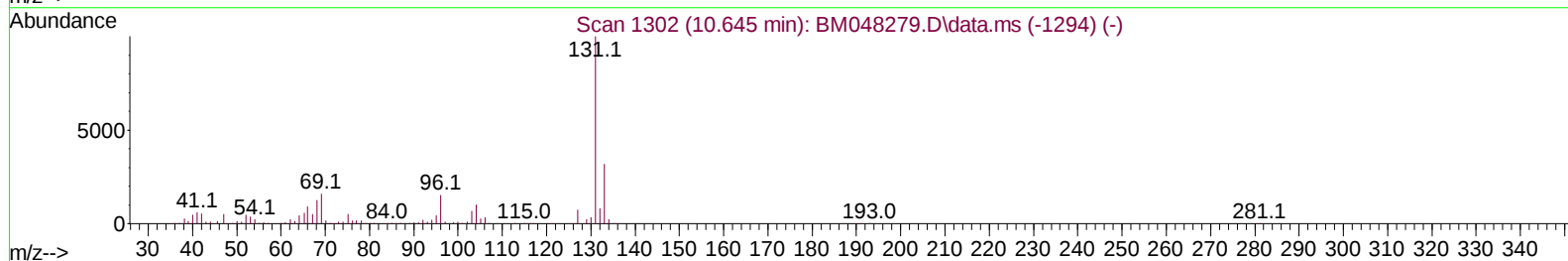
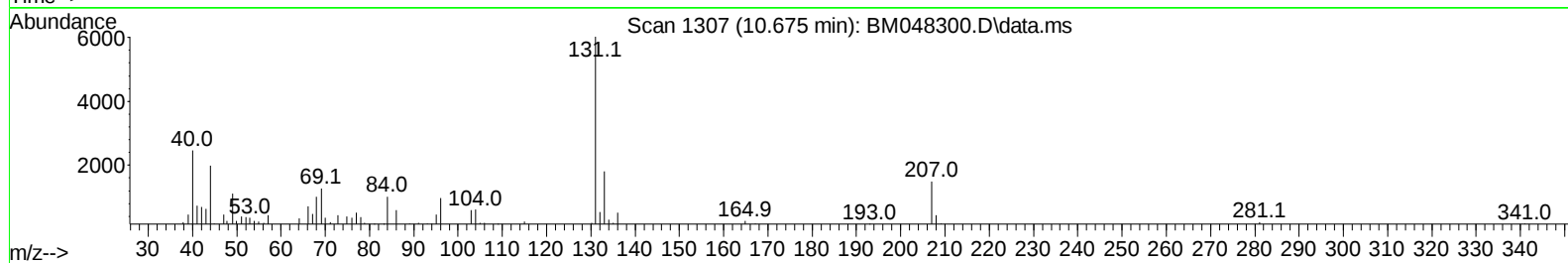
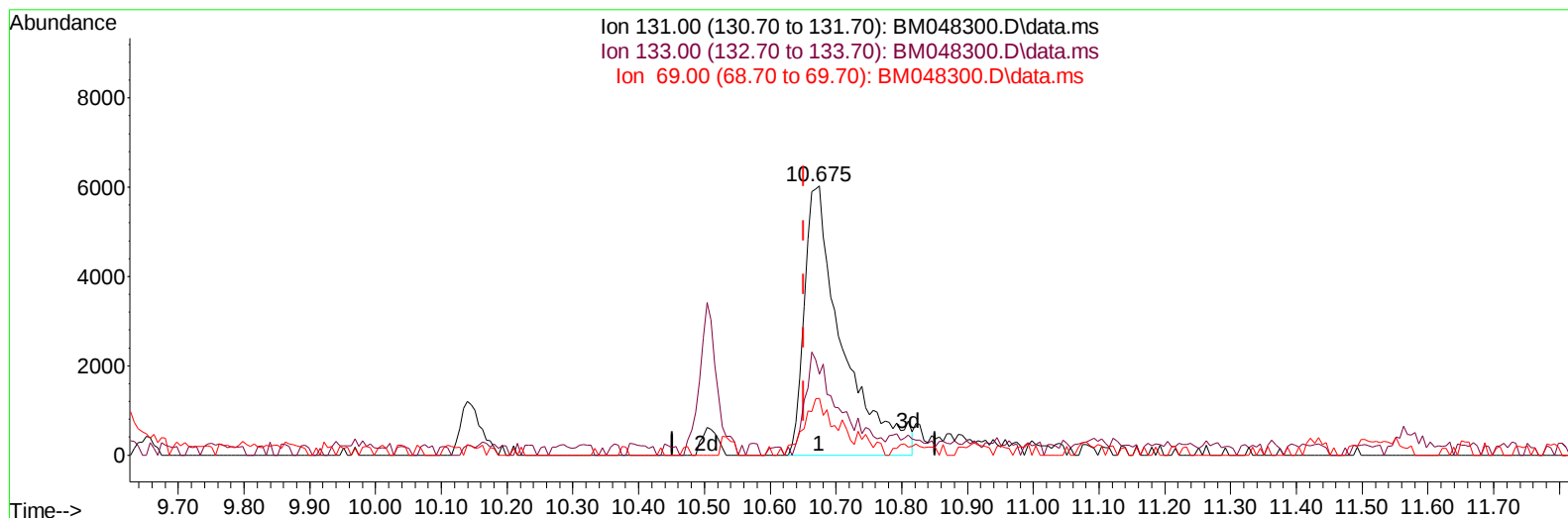
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM102824\  
Data File : BM048300.D  
Acq On : 29 Oct 2024 23:18  
Operator : RC/JU  
Sample : P4492-11  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
A4EQ6

Manual IntegrationsAPPROVED

Quant Time: Oct 29 23:51:42 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM102624.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun Oct 27 23:39:58 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/30/2024  
Supervised By :mohammad ahmed 11/01/2024



TIC: BM048300.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.675min (+ 0.024) 1.46 ng/ul m

response 24149

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 131.00 | 100.00 | 100.00 |
| 133.00 | 31.80  | 30.09  |
| 69.00  | 16.30  | 21.05# |
| 0.00   | 0.00   | 0.00   |



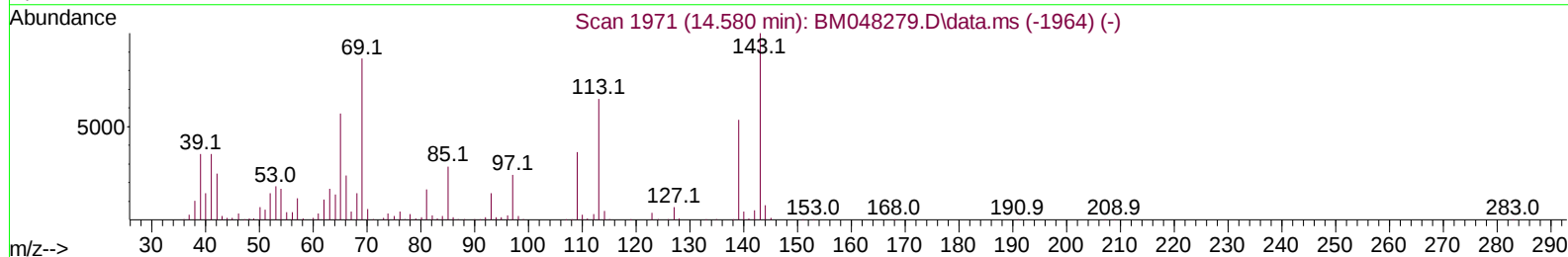
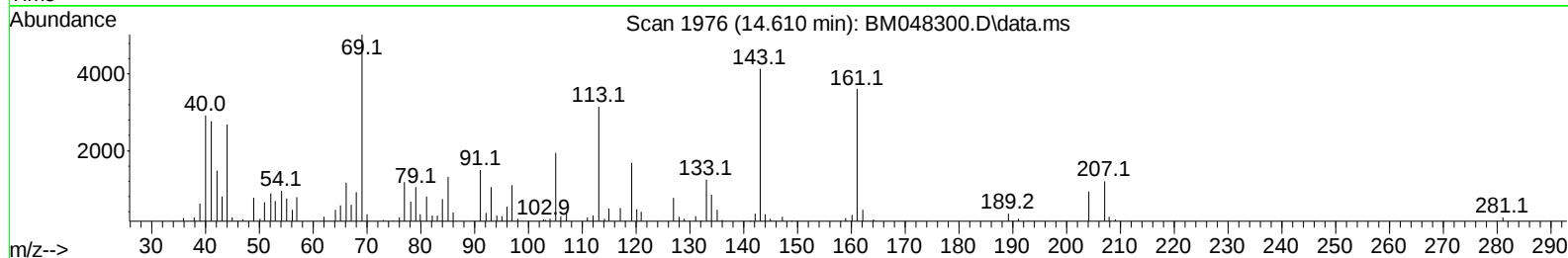
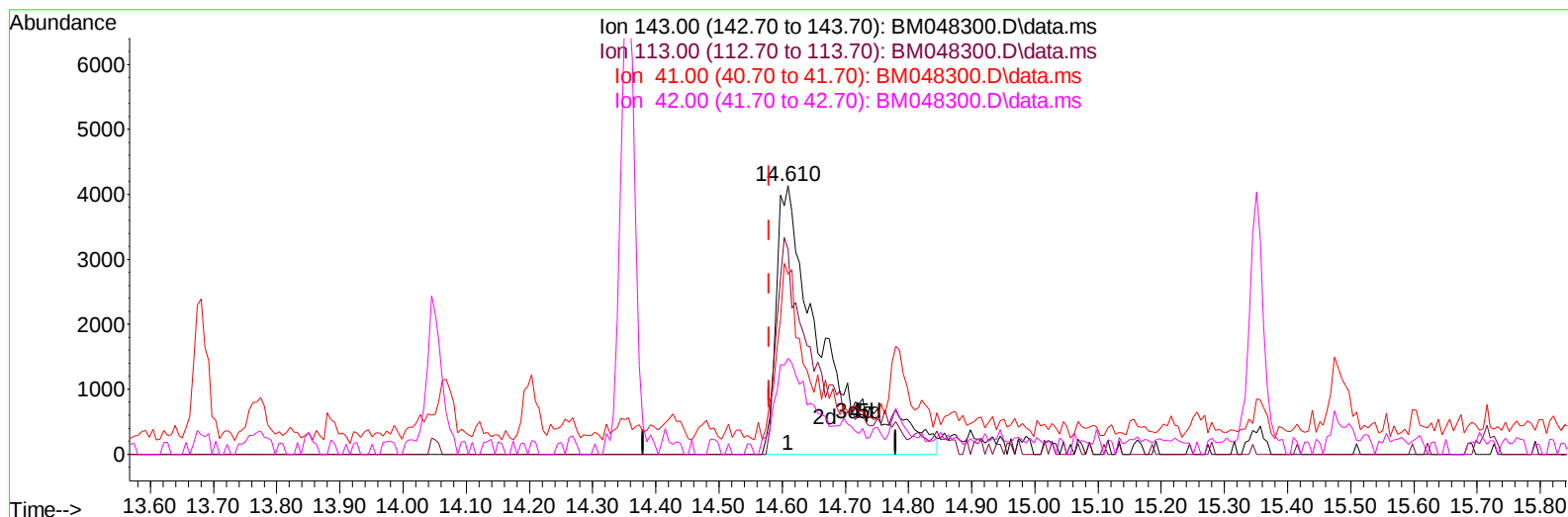
Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM102824\  
Data File : BM048300.D  
Acq On : 29 Oct 2024 23:18  
Operator : RC/JU  
Sample : P4492-11  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
A4EQ6

Manual IntegrationsAPPROVED

Quant Time: Oct 29 23:51:42 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM102624.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun Oct 27 23:39:58 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/30/2024  
Supervised By :mohammad ahmed 11/01/2024



TIC: BM048300.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.610min (+ 0.029) 3.58 ng/ul m

response 21325

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 143.00 | 100.00 | 100.00 |
| 113.00 | 68.30  | 76.31  |
| 41.00  | 40.60  | 67.06# |
| 42.00  | 26.40  | 35.77# |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM102824\  
 Data File : BM048300.D  
 Acq On : 29 Oct 2024 23:18  
 Operator : RC/JU  
 Sample : P4492-11  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 A4EQ6

Manual IntegrationsAPPROVED

Quant Time: Oct 29 23:53:44 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM102624.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Oct 27 23:39:58 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/30/2024  
 Supervised By :mohammad ahmed 11/01/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.716  | 152  | 180143   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.504 | 136  | 732971   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.357 | 164  | 435547   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.098 | 188  | 857802   | 20.000 | ng/ul  | -0.01    |
| 79) Chrysene-d12              | 21.321 | 240  | 790294   | 20.000 | ng/ul  | -0.01    |
| 88) Perylene-d12              | 24.262 | 264  | 871929   | 20.000 | ng/ul  | -0.02    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.169  | 96   | 6530     | 1.427  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/ul  |          |
| 7) Phenol-d5                  | 6.892  | 99   | 116386   | 7.520  | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.051  | 67   | 79921    | 8.337  | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.251  | 132  | 103812   | 8.248  | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.433  | 113  | 60071    | 4.835  | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.875  | 128  | 45797    | 7.621  | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.598  | 143  | 45594    | 7.004  | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.145 | 165  | 87196    | 7.340  | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.675 | 131  | 24149m   | 1.464  | ng/ul  | 0.02     |
| 46) Dimethylphthalate-d6      | 13.768 | 166  | 295230   | 9.070  | ng/ul  | -0.01    |
| 49) Acenaphthylene-d8         | 14.051 | 160  | 316503   | 8.624  | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.610 | 143  | 21325m   | 3.583  | ng/ul  | 0.03     |
| 60) Fluorene-d10              | 15.351 | 176  | 253778   | 9.227  | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.486 | 200  | 18704    | 3.299  | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.198 | 188  | 347241   | 8.561  | ng/ul  | -0.01    |
| 81) Pyrene-d10                | 19.492 | 212  | 408957   | 8.528  | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.062 | 264  | 315377   | 6.858  | ng/ul  | -0.02    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 8) Phenol                     | 6.922  | 94   | 23067    | 1.431  | ng/ul  | 95       |
| 16) Acetophenone              | 8.539  | 105  | 59746    | 3.053  | ng/ul  | 99       |
| 72) Phenanthrene              | 17.139 | 178  | 103581   | 2.181  | ng/ul  | 98       |
| 80) Fluoranthene              | 19.156 | 202  | 196405   | 3.520  | ng/ul  | 99       |
| 82) Pyrene                    | 19.521 | 202  | 148411   | 2.481  | ng/ul  | 97       |
| 85) Benzo(a)anthracene        | 21.303 | 228  | 82754    | 1.452  | ng/ul  | 96       |
| 87) Chrysene                  | 21.368 | 228  | 94469    | 1.763  | ng/ul  | 97       |
| 90) Benzo(b)fluoranthene      | 23.338 | 252  | 121883   | 2.190  | ng/ul# | 96       |
| -----                         |        |      |          |        |        |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM102824\  
Data File : BM048300.D  
Acq On : 29 Oct 2024 23:18  
Operator : RC/JU  
Sample : P4492-11  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
A4EQ6

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/30/2024  
Supervised By :mohammad ahmed 11/01/2024

Quant Time: Oct 29 23:53:44 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM102624.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun Oct 27 23:39:58 2024  
Response via : Initial Calibration

