

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP071923\  
 Data File : BP016321.D  
 Acq On : 19 Jul 2023 13:05  
 Operator : MA/JU  
 Sample : 03458-25DL 2X  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A46L5DL

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 07/19/2023  
 Supervised By :Sohil Jodhani 07/19/2023

Quant Time: Jul 19 15:04:04 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 19 01:28:59 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.958  | 152  | 95333m   | 20.000 | ng/u1  | 0.01     |
| 20) Naphthalene-d8            | 10.787 | 136  | 444702   | 20.000 | ng/u1  | 0.02     |
| 38) Acenaphthene-d10          | 14.610 | 164  | 303721   | 20.000 | ng/u1  | 0.01     |
| 64) Phenanthrene-d10          | 17.369 | 188  | 716533   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.457 | 240  | 853123   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.945 | 264  | 1033777  | 20.000 | ng/u1  | -0.02    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.281  | 96   | 3769m    | 1.399  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 7.222  | 99   | 35727    | 4.149  | ng/u1  | 0.09     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.316  | 67   | 31586m   | 5.425  | ng/u1  | 0.04     |
| 11) 2-Chlorophenol-d4         | 7.534  | 132  | 40555m   | 6.524  | ng/u1  | 0.06     |
| 15) 4-Methylphenol-d8         | 8.816  | 113  | 45905m   | 6.618  | ng/u1  | 0.12     |
| 21) Nitrobenzene-d5           | 9.216  | 128  | 19913m   | 6.491  | ng/u1  | 0.08     |
| 24) 2-Nitrophenol-d4          | 9.981  | 143  | 23587m   | 6.746  | ng/u1  | 0.12     |
| 28) 2,4-Dichlorophenol-d3     | 10.604 | 165  | 42102m   | 6.953  | ng/u1  | 0.18     |
| 31) 4-Chloroaniline-d4        | 11.263 | 131  | 5878m    | 0.606  | ng/u1  | 0.32     |
| 46) Dimethylphthalate-d6      | 14.051 | 166  | 181479   | 7.833  | ng/u1  | 0.04     |
| 49) Acenaphthylene-d8         | 14.310 | 160  | 165717   | 6.527  | ng/u1  | 0.01     |
| 54) 4-Nitrophenol-d4          | 15.069 | 143  | 18258m   | 3.861  | ng/u1  | 0.14     |
| 60) Fluorene-d10              | 15.622 | 176  | 159722   | 8.238  | ng/u1  | 0.02     |
| 65) 4,6-Dinitro-2-methylph... | 15.863 | 200  | 12733m   | 3.078  | ng/u1  | 0.08     |
| 73) Anthracene-d10            | 17.481 | 188  | 246132   | 7.853  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.704 | 212  | 177370   | 4.007  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.792 | 264  | 109915   | 2.214  | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 16) Acetophenone              | 8.893  | 105  | 39993m   | 3.496  | ng/u1  |          |
| 30) Naphthalene               | 10.846 | 128  | 32858    | 1.408  | ng/u1  | 98       |
| 52) Acenaphthene              | 14.681 | 153  | 76041    | 3.855  | ng/u1  | 100      |
| 56) Dibenzofuran              | 15.051 | 168  | 47892    | 1.783  | ng/u1  | 95       |
| 61) Fluorene                  | 15.687 | 166  | 76441    | 3.406  | ng/u1  | 97       |
| 72) Phenanthrene              | 17.410 | 178  | 1115764  | 29.886 | ng/u1  | 100      |
| 74) Anthracene                | 17.522 | 178  | 216200m  | 5.763  | ng/u1  |          |
| 77) Carbazole                 | 17.822 | 167  | 53875    | 1.575  | ng/u1  | 98       |
| 80) Fluoranthene              | 19.375 | 202  | 1895744  | 36.621 | ng/u1  | 97       |
| 82) Pyrene                    | 19.733 | 202  | 680142   | 12.306 | ng/u1  | 96       |
| 85) Benzo(a)anthracene        | 21.445 | 228  | 845567   | 15.220 | ng/u1  | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.322 | 149  | 48821    | 1.371  | ng/u1  | 100      |
| 87) Chrysene                  | 21.498 | 228  | 734185   | 13.046 | ng/u1  | 99       |
| 90) Benzo(b)fluoranthene      | 23.186 | 252  | 842456   | 14.267 | ng/u1  | 99       |
| 91) Benzo(k)fluoranthene      | 23.233 | 252  | 355896   | 5.922  | ng/u1  | 98       |
| 93) Benzo(a)pyrene            | 23.845 | 252  | 127327   | 2.177  | ng/u1  | 97       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.604 | 276  | 133771   | 1.847  | ng/u1  | 97       |
| 95) Dibenzo(a,h)anthracene    | 26.592 | 278  | 96106    | 1.627  | ng/u1  | 96       |
| -----                         |        |      |          |        |        |          |

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

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