

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM100324\  
 Data File : BM047842.D  
 Acq On : 04 Oct 2024 14:20  
 Operator : RC/JU  
 Sample : P4017-19ME  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DDCD5ME

Quant Time: Oct 04 14:54:46 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM092724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Oct 04 08:32:31 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.798  | 152  | 270126   | 20.000  | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.592 | 136  | 1093703  | 20.000  | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.433 | 164  | 696605   | 20.000  | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.174 | 188  | 1392752  | 20.000  | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.398 | 240  | 1321734  | 20.000  | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.397 | 264  | 1523817  | 20.000  | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 3.252  | 96   | 29434    | 3.521   | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.669  | 84   | 157745   | 6.445   | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.969  | 99   | 546492   | 21.395  | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 347226   | 19.157  | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.334  | 132  | 465797   | 25.064  | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.504  | 113  | 463790   | 24.421  | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.951  | 128  | 227158   | 26.086  | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.675  | 143  | 248597   | 29.790  | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.216 | 165  | 455535   | 27.326  | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.728 | 131  | 565899   | 21.958  | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.839 | 166  | 1444861  | 28.609  | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.127 | 160  | 1699076  | 28.374  | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.627 | 143  | 220889   | 21.967  | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.427 | 176  | 1286913  | 29.418  | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.539 | 200  | 173512   | 21.289  | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.268 | 188  | 1998468  | 29.182  | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.557 | 212  | 2314687  | 30.856  | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.191 | 264  | 2391101  | 30.243  | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |         |        |          |
|                               |        |      |          |         | Qvalue |          |
| 30) Naphthalene               | 10.639 | 128  | 88560    | 1.408   | ng/ul# | 96       |
| 36) 2-Methylnaphthalene       | 12.251 | 142  | 269906   | 6.752   | ng/ul  | 97       |
| 37) 1-Methylnaphthalene       | 12.475 | 142  | 149713   | 3.681   | ng/ul  | 97       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.057 | 232  | 411398   | 36.377  | ng/ul# | 97       |
| 71) Pentachlorophenol         | 16.833 | 266  | 3134810  | 295.084 | ng/ul  | 99       |
| 72) Phenanthrene              | 17.215 | 178  | 91632    | 1.122   | ng/ul  | 99       |
| -----                         |        |      |          |         |        |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM100324\  
Data File : BM047842.D  
Acq On : 04 Oct 2024 14:20  
Operator : RC/JU  
Sample : P4017-19ME  
Misc :  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
DDCD5ME

Quant Time: Oct 04 14:54:46 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM092724.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Oct 04 08:32:31 2024  
Response via : Initial Calibration

