

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN040424\  
 Data File : BN030643.D  
 Acq On : 04 Apr 2024 01:21  
 Operator : MA/JU  
 Sample : P1849-05  
 Misc :  
 ALS Vial : 32 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 EOLY9

Quant Time: Apr 04 01:55:04 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN032824.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Mar 31 12:24:07 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.693  | 152  | 241322   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.481 | 136  | 1064135  | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.340 | 164  | 665442   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.087 | 188  | 1227324  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.327 | 240  | 748909   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.280 | 264  | 800331   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.199  | 96   | 20778    | 4.037  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.605  | 84   | 91012    | 5.897  | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.864  | 99   | 437860   | 22.429 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.034  | 67   | 271474   | 22.515 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.228  | 132  | 371587   | 25.091 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.399  | 113  | 261260   | 15.968 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.846  | 128  | 193472   | 26.209 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.563  | 143  | 203572   | 30.145 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.105 | 165  | 375221   | 25.380 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.622 | 131  | 1643     | 0.068  | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.751 | 166  | 1156772  | 26.037 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.028 | 160  | 1354037  | 26.507 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.545 | 143  | 150539   | 17.714 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.334 | 176  | 1007379  | 25.911 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.451 | 200  | 71281    | 13.158 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.187 | 188  | 1350578  | 25.540 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.486 | 212  | 1365081  | 33.903 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.080 | 264  | 1009362  | 27.371 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 16) Acetophenone              | 8.516  | 105  | 78337    | 3.033  | ng/u1  | 99       |
| 47) Dimethylphthalate         | 13.798 | 163  | 444644   | 9.688  | ng/u1  | 99       |
| 72) Phenanthrene              | 17.128 | 178  | 105326   | 1.670  | ng/u1  | 99       |
| 80) Fluoranthene              | 19.151 | 202  | 239481   | 5.003  | ng/u1  | 99       |
| 82) Pyrene                    | 19.516 | 202  | 193917   | 3.836  | ng/u1  | 99       |
| 83) Butylbenzylphthalate      | 20.428 | 149  | 40966    | 2.122  | ng/u1# | 96       |
| 85) Benzo(a)anthracene        | 21.310 | 228  | 78701    | 1.575  | ng/u1# | 89       |
| 86) Bis(2-ethylhexyl)phtha... | 21.245 | 149  | 215973   | 7.586  | ng/u1  | 99       |
| 87) Chrysene                  | 21.369 | 228  | 115964   | 2.458  | ng/u1  | 96       |
| 90) Benzo(b)fluoranthene      | 23.345 | 252  | 161894   | 3.437  | ng/u1# | 74       |
| 91) Benzo(k)fluoranthene      | 23.404 | 252  | 48143    | 1.006  | ng/u1# | 63       |
| 93) Benzo(a)pyrene            | 24.133 | 252  | 75448    | 1.696  | ng/u1# | 74       |
| 94) Indeno(1,2,3-cd)pyrene    | 27.592 | 276  | 83942    | 1.745  | ng/u1# | 87       |
| 96) Benzo(g,h,i)perylene      | 28.621 | 276  | 91650    | 2.427  | ng/u1  | 92       |

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

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