

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012725\  
 Data File : BP023743.D  
 Acq On : 27 Jan 2025 16:31  
 Operator : RC/JU  
 Sample : Q1174-07  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 E2937

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 01/28/2025  
 Supervised By :mohammad ahmed 01/31/2025

Quant Time: Jan 27 23:10:45 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012225.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 27 21:53:59 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.793  | 152  | 655120   | 20.000  | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.569 | 136  | 2802119  | 20.000  | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.428 | 164  | 1894862  | 20.000  | ng/u1  | 0.01     |
| 64) Phenanthrene-d10          | 17.228 | 188  | 3592714  | 20.000  | ng/u1  | 0.01     |
| 79) Chrysene-d12              | 21.680 | 240  | 3599804  | 20.000  | ng/u1  | 0.00     |
| 88) Perylene-d12              | 25.092 | 264  | 4004453  | 20.000  | ng/u1  | 0.03     |
|                               |        |      |          |         |        |          |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 72112    | 4.433   | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.716  | 84   | 508506   | 10.834  | ng/u1  | 0.01     |
| 7) Phenol-d5                  | 6.975  | 99   | 628944   | 11.346  | ng/u1  | 0.02     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 1026092  | 32.771  | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.334  | 132  | 1487157  | 35.289  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.499  | 113  | 1132765  | 25.987  | ng/u1  | 0.02     |
| 21) Nitrobenzene-d5           | 8.940  | 128  | 821896   | 38.930  | ng/u1  | 0.01     |
| 24) 2-Nitrophenol-d4          | 9.657  | 143  | 920957   | 47.582  | ng/u1  | 0.01     |
| 28) 2,4-Dichlorophenol-d3     | 10.199 | 165  | 1634073  | 40.165  | ng/u1  | 0.02     |
| 31) 4-Chloroaniline-d4        | 10.704 | 131  | 1327617  | 20.405  | ng/u1  | 0.01     |
| 46) Dimethylphthalate-d6      | 13.834 | 166  | 4712844  | 34.584  | ng/u1  | 0.02     |
| 49) Acenaphthylene-d8         | 14.116 | 160  | 5677215  | 35.996  | ng/u1  | 0.01     |
| 54) 4-Nitrophenol-d4          | 14.645 | 143  | 301636   | 11.016  | ng/u1  | 0.05     |
| 60) Fluorene-d10              | 15.439 | 176  | 4344582  | 37.244  | ng/u1  | 0.02     |
| 65) 4,6-Dinitro-2-methylph... | 15.551 | 200  | 757842   | 42.835  | ng/u1  | 0.02     |
| 73) Anthracene-d10            | 17.333 | 188  | 6695965  | 38.821  | ng/u1  | 0.01     |
| 81) Pyrene-d10                | 19.698 | 212  | 7453409  | 40.690  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.868 | 264  | 7791439  | 38.873  | ng/u1  | 0.02     |
|                               |        |      |          |         |        |          |
| Target Compounds              |        |      |          |         |        |          |
| 2) 1,4-Dioxane                | 3.334  | 88   | 72909    | 4.237   | ng/uL# | 95       |
| 8) Phenol                     | 7.010  | 94   | 10661607 | 183.807 | ng/u1  | 97       |
| 13) 2-Methylphenol            | 8.240  | 108  | 59906    | 1.423   | ng/u1  | 100      |
| 18) 4-Methylphenol            | 8.563  | 108  | 1470059  | 31.411  | ng/u1  | 98       |
| 26) 2,4-Dimethylphenol        | 9.787  | 107  | 682892m  | 14.499  | ng/u1  |          |
| 36) 2-Methylnaphthalene       | 12.234 | 142  | 170636   | 1.676   | ng/u1  | 96       |
| 37) 1-Methylnaphthalene       | 12.451 | 142  | 156268   | 1.535   | ng/u1  | 99       |
| 42) 2,4,5-Trichlorophenol     | 12.922 | 196  | 55838    | 1.528   | ng/u1  | 96       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.063 | 232  | 167336   | 5.575   | ng/u1# | 89       |
| 71) Pentachlorophenol         | 16.875 | 266  | 29672    | 1.049   | ng/u1  | 97       |
| -----                         |        |      |          |         |        |          |

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

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