

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN021424\  
 Data File : BN029794.D  
 Acq On : 15 Feb 2024 06:28  
 Operator : MA/JU  
 Sample : P1329-06  
 Misc :  
 ALS Vial : 73 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 C0FL5

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 02/16/2024  
 Supervised By :mohammad ahmed 02/17/2024

Quant Time: Feb 15 06:55:31 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN020724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 13 22:50:45 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.869  | 152  | 169605   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.663 | 136  | 688198   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.510 | 164  | 370563   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.263 | 188  | 766967   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.463 | 240  | 688035   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.839 | 264  | 765035   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.323  | 96   | 18902    | 3.666  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.734  | 84   | 165008   | 11.999 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.040  | 99   | 345723   | 21.793 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.211  | 67   | 221619   | 20.932 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.399  | 132  | 273491   | 23.790 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.581  | 113  | 114250   | 9.657  | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.034  | 128  | 136417   | 25.540 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.752  | 143  | 148001   | 31.815 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.287 | 165  | 230890   | 25.146 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.816 | 131  | 282317   | 18.022 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.934 | 166  | 777509   | 28.182 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.204 | 160  | 890351   | 27.175 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.722 | 143  | 121751   | 23.251 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.510 | 176  | 662988   | 28.248 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.634 | 200  | 77695    | 20.790 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.363 | 188  | 954324   | 26.550 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.663 | 212  | 1291724  | 31.285 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.686 | 264  | 1208277  | 31.563 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 16) Acetophenone              | 8.699  | 105  | 39313    | 2.088  | ng/u1  | 98       |
| 72) Phenanthrene              | 17.304 | 178  | 276215   | 6.226  | ng/u1  | 99       |
| 74) Anthracene                | 17.398 | 178  | 51039    | 1.146  | ng/u1  | 95       |
| 80) Fluoranthene              | 19.328 | 202  | 728727   | 14.529 | ng/u1  | 98       |
| 82) Pyrene                    | 19.692 | 202  | 646040   | 12.176 | ng/u1  | 95       |
| 85) Benzo(a)anthracene        | 21.451 | 228  | 357286   | 7.069  | ng/u1  | 97       |
| 87) Chrysene                  | 21.504 | 228  | 330360   | 6.974  | ng/u1  | 96       |
| 90) Benzo(b)fluoranthene      | 23.116 | 252  | 392271   | 8.141  | ng/u1  | 96       |
| 91) Benzo(k)fluoranthene      | 23.163 | 252  | 163463m  | 3.314  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 23.733 | 252  | 255188   | 5.522  | ng/u1  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.280 | 276  | 145283   | 2.582  | ng/u1  | 97       |
| 96) Benzo(g,h,i)perylene      | 27.027 | 276  | 120970   | 2.570  | ng/u1  | 95       |

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

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