

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102521\  
 Data File : BP007665.D  
 Acq On : 26 Oct 2021 05:00  
 Operator : CG/JU  
 Sample : M4291-10  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 P022-TW001-02

Manual Integrations  
 APPROVED

mohammad  
 10/26/2021 11:35:00 AM

Quant Time: Oct 26 05:49:34 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP102521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Oct 25 16:57:16 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.934  | 152  | 172697   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.740 | 136  | 694229   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.575 | 164  | 460866   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.328 | 188  | 1026975  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.416 | 240  | 1133911  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.857 | 264  | 1183430  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.370  | 96   | 21166    | 5.567  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.793  | 84   | 52051    | 4.527  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.099  | 99   | 109561   | 8.352  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.264  | 67   | 249749   | 34.293 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.464  | 132  | 284311   | 27.868 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.646  | 113  | 191066   | 18.602 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.093  | 128  | 159934   | 33.806 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.816  | 143  | 164052   | 31.749 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.358 | 165  | 308177   | 29.301 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.875 | 131  | 348087   | 25.030 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.981 | 166  | 1113656  | 35.674 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.263 | 160  | 1276782  | 33.369 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.763 | 143  | 41078m   | 7.178  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.569 | 176  | 964799   | 36.660 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.681 | 200  | 151528   | 27.584 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.428 | 188  | 1663453  | 39.414 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.657 | 212  | 2062563  | 39.070 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.698 | 264  | 2139501  | 36.802 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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