

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP030921\  
 Data File : BP004962.D  
 Acq On : 10 Mar 2021 10:31  
 Operator : CG/JU  
 Sample : PB134967BL  
 Misc : SOM-MDL-W-BL-07  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK67

Quant Time: Mar 10 11:03:49 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP030921MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Mar 10 10:08:47 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.74  | 152  | 31791    | 20.00 | ng/ul | 0.01     |
| 18) Naphthalene-d8        | 10.53 | 136  | 127565   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.39 | 164  | 84207    | 20.00 | ng/ul | 0.01     |
| 62) Phenanthrene-d10      | 17.15 | 188  | 190605   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.23 | 240  | 209212   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.53 | 264  | 217843   | 20.00 | ng/ul | 0.01     |

| System Monitoring Compounds    | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 3) 1,4-Dioxane-d8              | 3.23  | 96   | 5666     | 6.66  | ng/uL | 0.01     |
| 5) Phenol-d5                   | 6.91  | 99   | 66696    | 26.71 | ng/ul | 0.00     |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.08  | 67   | 38016    | 28.66 | ng/ul | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.27  | 132  | 52647    | 26.69 | ng/ul | 0.00     |
| 13) 4-Methylphenol-d8          | 8.45  | 113  | 54509    | 25.91 | ng/ul | 0.00     |
| 19) Nitrobenzene-d5            | 8.89  | 128  | 26042    | 27.72 | ng/ul | 0.00     |
| 22) 2-Nitrophenol-d4           | 9.62  | 143  | 29512    | 27.99 | ng/ul | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 10.15 | 165  | 50146    | 24.72 | ng/ul | 0.00     |
| 29) 4-Chloroaniline-d4         | 10.67 | 131  | 71121    | 32.03 | ng/ul | 0.00     |
| 44) Dimethylphthalate-d6       | 13.80 | 166  | 204874   | 33.06 | ng/ul | 0.01     |
| 47) Acenaphthylene-d8          | 14.08 | 160  | 218800   | 28.93 | ng/ul | 0.00     |
| 52) 4-Nitrophenol-d4           | 14.60 | 143  | 27642    | 25.96 | ng/ul | 0.00     |
| 58) Fluorene-d10               | 15.39 | 176  | 165053   | 30.98 | ng/ul | 0.00     |
| 63) 4,6-Dinitro-2-methylphenol | 15.51 | 200  | 34903    | 28.55 | ng/ul | 0.01     |
| 71) Anthracene-d10             | 17.25 | 188  | 286621   | 33.55 | ng/ul | 0.00     |
| 79) Pyrene-d10                 | 19.49 | 212  | 343085   | 35.36 | ng/ul | 0.00     |
| 90) Benzo(a)pyrene-d12         | 23.38 | 264  | 383152   | 34.39 | ng/ul | 0.01     |

| Target Compounds | Ovalue |
|------------------|--------|
| -----            |        |

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

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