

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP031220\  
 Data File : BP002201.D  
 Acq On : 13 Mar 2020 00:57  
 Operator : CG/JU  
 Sample : TESTBL-03  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TESTBL-03

Manual Integrations  
 APPROVED

mohammad  
 3/13/2020 8:57:34 AM

Quant Time: Mar 13 04:06:20 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP022720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 13 02:52:52 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.63  | 152  | 271464   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.40 | 136  | 1121892  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.26 | 164  | 696526   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.02 | 188  | 1493969  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.11 | 240  | 1462654  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.31 | 264  | 1695551  | 20.00 | ng/ul | 0.00     |

| System Monitoring Compounds    | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 3) 1,4-Dioxane-d8              | 3.15  | 96   | 37357    | 5.92  | ng/uL | 0.00     |
| 5) Phenol-d5                   | 6.81  | 99   | 639328   | 30.87 | ng/ul | 0.00     |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.96  | 67   | 372492   | 33.97 | ng/ul | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.16  | 132  | 565803   | 31.58 | ng/ul | 0.00     |
| 13) 4-Methylphenol-d8          | 8.33  | 113  | 515225   | 30.28 | ng/ul | 0.00     |
| 19) Nitrobenzene-d5            | 8.77  | 128  | 252438   | 29.29 | ng/ul | 0.00     |
| 22) 2-Nitrophenol-d4           | 9.49  | 143  | 272055   | 28.06 | ng/ul | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 10.03 | 165  | 525977   | 28.16 | ng/ul | 0.00     |
| 29) 4-Chloroaniline-d4         | 10.53 | 131  | 726531   | 35.48 | ng/ul | 0.00     |
| 44) Dimethylphthalate-d6       | 13.68 | 166  | 1572764  | 29.84 | ng/ul | 0.00     |
| 47) Acenaphthylene-d8          | 13.95 | 160  | 1940178  | 30.78 | ng/ul | 0.00     |
| 52) 4-Nitrophenol-d4           | 14.47 | 143  | 198565   | 19.70 | ng/ul | 0.00     |
| 58) Fluorene-d10               | 15.26 | 176  | 1443475  | 31.65 | ng/ul | 0.00     |
| 63) 4,6-Dinitro-2-methylphenol | 15.39 | 200  | 163916   | 19.90 | ng/ul | 0.00     |
| 71) Anthracene-d10             | 17.11 | 188  | 2254376m | 33.57 | ng/ul | 0.00     |
| 79) Pyrene-d10                 | 19.36 | 212  | 2590804  | 33.84 | ng/ul | 0.00     |
| 90) Benzo(a)pyrene-d12         | 23.17 | 264  | 2875425  | 33.15 | ng/ul | 0.00     |

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

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

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