

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
 Data File : BP002790.D  
 Acq On : 17 Jul 2020 14:03  
 Operator : CG/JU  
 Sample : L3324-01  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 10:1-COMP

Quant Time: Jul 17 14:52:58 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jul 13 10:54:03 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.56  | 152  | 230785   | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 10.33 | 136  | 943594   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 14.21 | 164  | 454315   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 16.97 | 188  | 658129   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 21.09 | 240  | 565896   | 20.00 | ng    | 0.01     |
| 86) Perylene-d12          | 23.27 | 264  | 642354   | 20.00 | ng    | 0.01     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|-------|------|----------|-------|-------|----------|
| 5) 2-Fluorophenol           | 5.19  | 112  | 1100420  | 80.45 | ng    | 0.00     |
| 7) Phenol-d6                | 6.76  | 99   | 1497053  | 76.69 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.70  | 82   | 965711   | 54.74 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.72 | 330  | 337967   | 71.27 | ng    | 0.01     |
| 45) 2-Fluorobiphenyl        | 12.82 | 172  | 1774456  | 60.42 | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.57 | 244  | 1142684  | 46.30 | ng    | 0.01     |

| Target Compounds               | R.T.  | QIon | Response | Conc  | Units | Qvalue |
|--------------------------------|-------|------|----------|-------|-------|--------|
| 50) Dimethylphthalate          | 13.67 | 163  | 334586   | 9.657 | ng    | # 98   |
| 71) Phenanthrene               | 17.02 | 178  | 99652    | 2.797 | ng    | # 69   |
| 75) Fluoranthene               | 19.01 | 202  | 248639   | 6.022 | ng    | 98     |
| 78) Pyrene                     | 19.36 | 202  | 256460   | 6.564 | ng    | 98     |
| 81) Benzo(a)anthracene         | 21.07 | 228  | 129001   | 3.522 | ng    | 94     |
| 83) Chrysene                   | 21.12 | 228  | 174803   | 5.013 | ng    | 97     |
| 84) Bis(2-ethylhexyl)phthalate | 21.01 | 149  | 74434    | 3.471 | ng    | 92     |
| 87) Indeno(1,2,3-cd)pyrene     | 25.48 | 276  | 118201   | 2.546 | ng    | 96     |
| 88) Benzo(b)fluoranthene       | 22.62 | 252  | 261167m  | 6.548 | ng    |        |
| 90) Benzo(a)pyrene             | 23.18 | 252  | 149439   | 3.954 | ng    | # 92   |
| 92) Benzo(g,h,i)perylene       | 26.16 | 276  | 117447   | 3.094 | ng    | 95     |

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

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