

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080719\  
 Data File : BF116051.D  
 Acq On : 7 Aug 2019 17:07  
 Operator : HP/JU  
 Sample : K4152-11  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 2S-E1-E4-COMP-(0-1)

Manual Integrations  
 APPROVED

mohammad  
 8/8/2019 4:46:53 PM

Quant Time: Aug 08 08:02:13 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Aug 07 11:13:18 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.85  | 152  | 142119   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.12  | 136  | 558028   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.87  | 164  | 291441   | 20.00  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.36 | 188  | 471318   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.00 | 240  | 266009   | 20.00  | ng    | 0.00     |
| 87) Perylene-d12            | 15.47 | 264  | 331182   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.47  | 112  | 964109   | 113.57 | ng    | 0.00     |
| 7) Phenol-d6                | 6.48  | 99   | 1260289  | 117.66 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.40  | 82   | 827053   | 85.55  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.67 | 330  | 263347   | 93.20  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.20  | 172  | 1278559  | 82.17  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.95 | 244  | 1101788  | 87.28  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       |          |
| 50) Dimethylphthalate       | 9.59  | 163  | 293375   | 14.784 | ng    | 99       |
| 71) Phenanthrene            | 11.39 | 178  | 301659   | 12.520 | ng    | 99       |
| 72) Anthracene              | 11.44 | 178  | 97719    | 4.007  | ng    | 97       |
| 75) Fluoranthene            | 12.58 | 202  | 424925   | 16.845 | ng    | 88       |
| 78) Pyrene                  | 12.81 | 202  | 342674   | 17.089 | ng    | # 88     |
| 81) Benzo(a)anthracene      | 14.00 | 228  | 152613   | 8.976  | ng    | 97       |
| 83) Chrysene                | 14.03 | 228  | 134155   | 8.127  | ng    | 94       |
| 86) Indeno(1,2,3-cd)pyrene  | 16.94 | 276  | 80028    | 5.772  | ng    | 95       |
| 88) Benzo(b)fluoranthene    | 15.04 | 252  | 167299m  | 8.737  | ng    |          |
| 89) Benzo(k)fluoranthene    | 15.07 | 252  | 58834m   | 3.154  | ng    |          |
| 90) Benzo(a)pyrene          | 15.41 | 252  | 126778   | 7.406  | ng    | 100      |
| 92) Benzo(a,h,i)perylene    | 17.38 | 276  | 77476    | 5.324  | ng    | 97       |

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

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