

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030220\  
 Data File : BF119175.D  
 Acq On : 2 Mar 2020 18:22  
 Operator : CG/JU  
 Sample : L1372-06 2X  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SU-03-022820

Manual Integrations  
 APPROVED

mohammad  
 3/3/2020 12:25:56 PM

Quant Time: Mar 03 08:04:55 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Feb 26 16:03:05 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.73  | 152  | 140038   | 20.00  | ng    | -0.01    |
| 21) Naphthalene-d8          | 8.02  | 136  | 548689   | 20.00  | ng    | -0.02    |
| 39) Acenaphthene-d10        | 9.77  | 164  | 294846   | 20.00  | ng    | -0.02    |
| 64) Phenanthrene-d10        | 11.25 | 188  | 500910   | 20.00  | ng    | -0.01    |
| 76) Chrysene-d12            | 13.89 | 240  | 348111   | 20.00  | ng    | -0.01    |
| 87) Perylene-d12            | 15.32 | 264  | 348495   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.36  | 112  | 363180   | 43.34  | ng    | -0.01    |
| 7) Phenol-d6                | 6.38  | 99   | 451950   | 44.35  | ng    | -0.02    |
| 23) Nitrobenzene-d5         | 7.29  | 82   | 294468   | 28.34  | ng    | -0.03    |
| 42) 2,4,6-Tribromophenol    | 10.56 | 330  | 144171   | 37.38  | ng    | -0.02    |
| 45) 2-Fluorobiphenyl        | 9.09  | 172  | 604324   | 30.19  | ng    | -0.02    |
| 79) Terphenyl-d14           | 12.83 | 244  | 630378   | 31.18  | ng    | -0.02    |
| Target Compounds            |       |      |          |        |       |          |
| 49) Acenaphthylene          | 9.63  | 152  | 65814    | 2.373  | ng    | 99       |
| 71) Phenanthrene            | 11.27 | 178  | 187475   | 6.863  | ng    | 96       |
| 72) Anthracene              | 11.33 | 178  | 69890    | 2.561  | ng    | 97       |
| 75) Fluoranthene            | 12.46 | 202  | 509355   | 17.566 | ng    | 99       |
| 78) Pyrene                  | 12.69 | 202  | 486537   | 17.406 | ng    | 98       |
| 81) Benzo(a)anthracene      | 13.88 | 228  | 338914   | 14.289 | ng    | 97       |
| 83) Chrysene                | 13.92 | 228  | 268536   | 11.362 | ng    | 98       |
| 86) Indeno(1,2,3-cd)pyrene  | 16.72 | 276  | 154232   | 6.740  | ng    | 96       |
| 88) Benzo(b)fluoranthene    | 14.91 | 252  | 453439m  | 19.740 | ng    |          |
| 89) Benzo(k)fluoranthene    | 14.93 | 252  | 149826m  | 7.038  | ng    |          |
| 90) Benzo(a)pyrene          | 15.26 | 252  | 303776   | 14.653 | ng    | 98       |
| 91) Dibenzo(a,h)anthracene  | 16.72 | 278  | 42759    | 2.344  | ng    | # 78     |
| 92) Benzo(a,h,i)perylene    | 17.15 | 276  | 136099   | 7.551  | ng    | 94       |

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

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