

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080719\  
 Data File : BF116043.D  
 Acq On : 7 Aug 2019 11:54  
 Operator : HP/JU  
 Sample : K4152-06  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 2S-H1-H4-COMP-(30)

Manual Integrations  
 APPROVED

mohammad  
 8/8/2019 4:46:43 PM

Quant Time: Aug 07 14:22:15 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   | 6.84  | 152  | 140523   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.12  | 136  | 541882   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.87  | 164  | 292223   | 20.00  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.36 | 188  | 508880   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.00 | 240  | 311475   | 20.00  | ng    | 0.00     |
| 87) Perylene-d12            | 15.47 | 264  | 294871   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.47  | 112  | 896532   | 106.80 | ng    | 0.00     |
| 7) Phenol-d6                | 6.48  | 99   | 1154744  | 109.03 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.40  | 82   | 739078   | 78.73  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.67 | 330  | 197626   | 69.75  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.20  | 172  | 1191477  | 76.37  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.95 | 244  | 1261995  | 85.38  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       |          |
| 50) Dimethylphthalate       | 9.59  | 163  | 288225   | 14.486 | ng    | 98       |
| 71) Phenanthrene            | 11.39 | 178  | 132700   | 5.101  | ng    | 98       |
| 74) Di-n-butylphthalate     | 11.96 | 149  | 137354   | 4.929  | ng    | # 97     |
| 75) Fluoranthene            | 12.58 | 202  | 219549   | 8.061  | ng    | 92       |
| 78) Pyrene                  | 12.81 | 202  | 183152   | 7.801  | ng    | # 94     |
| 81) Benzo(a)anthracene      | 13.99 | 228  | 84671    | 4.253  | ng    | 97       |
| 83) Chrysene                | 14.03 | 228  | 65585    | 3.393  | ng    | 96       |
| 86) Indeno(1,2,3-cd)pyrene  | 16.94 | 276  | 33990    | 2.094  | ng    | # 92     |
| 88) Benzo(b)fluoranthene    | 15.04 | 252  | 77684m   | 4.556  | ng    |          |
| 90) Benzo(a)pyrene          | 15.40 | 252  | 57083    | 3.745  | ng    | # 94     |
| 92) Benzo(g,h,i)perylene    | 17.39 | 276  | 32825    | 2.533  | ng    | # 94     |

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

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