

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112719\  
 Data File : BF118038.D  
 Acq On : 27 Nov 2019 14:04  
 Operator : JU  
 Sample : K5973-04  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 DUCT-BANK-2

Manual Integrations  
 APPROVED

mohammad  
 11/29/2019 8:53:17 AM

Quant Time: Nov 27 14:39:53 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 26 05:38:20 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4      | 6.88  | 152  | 142194   | 20.00  | ng    | -0.01    |
| 21) Naphthalene-d8             | 8.17  | 136  | 547901   | 20.00  | ng    | -0.01    |
| 39) Acenaphthene-d10           | 9.93  | 164  | 283261   | 20.00  | ng    | -0.01    |
| 64) Phenanthrene-d10           | 11.42 | 188  | 505021   | 20.00  | ng    | -0.01    |
| 76) Chrysene-d12               | 14.07 | 240  | 333361   | 20.00  | ng    | -0.01    |
| 87) Perylene-d12               | 15.57 | 264  | 393206   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds    |       |      |          |        |       |          |
| 5) 2-Fluorophenol              | 5.49  | 112  | 680109   | 84.66  | ng    | 0.00     |
| 7) Phenol-d6                   | 6.51  | 99   | 847602   | 87.48  | ng    | -0.01    |
| 23) Nitrobenzene-d5            | 7.44  | 82   | 520609   | 56.48  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol       | 10.72 | 330  | 260952   | 87.02  | ng    | -0.02    |
| 45) 2-Fluorobiphenyl           | 9.25  | 172  | 1031549  | 65.33  | ng    | -0.02    |
| 79) Terphenyl-d14              | 13.02 | 244  | 1061506  | 58.20  | ng    | -0.01    |
| Target Compounds               |       |      |          |        |       | Qvalue   |
| 50) Dimethylphthalate          | 9.64  | 163  | 213976   | 10.772 | ng    | 98       |
| 52) Acenaphthene               | 9.96  | 154  | 41126    | 2.547  | ng    | 97       |
| 58) Fluorene                   | 10.48 | 166  | 42636    | 2.461  | ng    | 99       |
| 71) Phenanthrene               | 11.45 | 178  | 830057   | 32.691 | ng    | 98       |
| 72) Anthracene                 | 11.50 | 178  | 132778   | 5.274  | ng    | 98       |
| 73) Carbazole                  | 11.65 | 167  | 78838    | 3.584  | ng    | 95       |
| 75) Fluoranthene               | 12.65 | 202  | 1063946  | 46.219 | ng    | 99       |
| 78) Pyrene                     | 12.87 | 202  | 943884   | 35.036 | ng    | 99       |
| 81) Benzo(a)anthracene         | 14.06 | 228  | 459991   | 20.881 | ng    | 100      |
| 83) Chrysene                   | 14.10 | 228  | 446111   | 20.899 | ng    | 98       |
| 84) Bis(2-ethylhexyl)phthalate | 14.04 | 149  | 41263    | 2.941  | ng    | 93       |
| 86) Indeno(1,2,3-cd)pyrene     | 17.09 | 276  | 286530m  | 15.771 | ng    |          |
| 88) Benzo(b)fluoranthene       | 15.13 | 252  | 584526m  | 22.881 | ng    |          |
| 89) Benzo(k)fluoranthene       | 15.16 | 252  | 167746m  | 6.969  | ng    |          |
| 90) Benzo(a)pyrene             | 15.51 | 252  | 407015   | 18.118 | ng    | 98       |
| 91) Dibenzo(a,h)anthracene     | 17.10 | 278  | 74060    | 3.830  | ng    | 96       |
| 92) Benzo(a,h,i)perylene       | 17.54 | 276  | 272220   | 14.135 | ng    | 96       |

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

