

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\  
 Data File : BP001240.D  
 Acq On : 06 Dec 2019 20:22  
 Operator : JU  
 Sample : K6112-01  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 OK-02-120419

Manual Integrations  
 APPROVED

mohammad  
 12/9/2019 3:10:53 PM

Quant Time: Dec 07 02:45:44 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Dec 05 12:34:47 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min)  |
|-----------------------------|-------|------|----------|--------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4   | 8.31  | 152  | 79640    | 20.00  | ng    | 0.00      |
| 21) Naphthalene-d8          | 10.34 | 136  | 293242   | 20.00  | ng    | -0.01     |
| 39) Acenaphthene-d10        | 13.20 | 164  | 171852   | 20.00  | ng    | -0.01     |
| 64) Phenanthrene-d10        | 15.60 | 188  | 322207   | 20.00  | ng    | -0.01     |
| 76) Chrysene-d12            | 19.24 | 240  | 239027   | 20.00  | ng    | -0.02     |
| 87) Perylene-d12            | 21.02 | 264  | 255675   | 20.00  | ng    | -0.01     |
| System Monitoring Compounds |       |      |          |        |       |           |
| 5) 2-Fluorophenol           | 6.22  | 112  | 446646   | 93.05  | ng    | 0.00      |
| 7) Phenol-d6                | 7.76  | 99   | 621602   | 97.20  | ng    | 0.00      |
| 23) Nitrobenzene-d5         | 9.19  | 82   | 424944   | 62.87  | ng    | -0.01     |
| 42) 2,4,6-Tribromophenol    | 14.49 | 330  | 168270   | 102.16 | ng    | -0.01     |
| 45) 2-Fluorobiphenyl        | 12.12 | 172  | 721795   | 61.47  | ng    | -0.01     |
| 79) Terphenyl-d14           | 17.88 | 244  | 811698   | 62.83  | ng    | -0.01     |
| Target Compounds            |       |      |          |        |       |           |
| 49) Acenaphthylene          | 12.96 | 152  | 33948    | 2.135  | ng    | Qvalue 97 |
| 50) Dimethylphthalate       | 12.79 | 163  | 106251   | 8.267  | ng    | 98        |
| 71) Phenanthrene            | 15.63 | 178  | 58298    | 3.442  | ng    | 97        |
| 75) Fluoranthene            | 17.34 | 202  | 111456   | 6.215  | ng    | 98        |
| 78) Pyrene                  | 17.65 | 202  | 143149   | 7.604  | ng    | 97        |
| 81) Benzo(a)anthracene      | 19.23 | 228  | 73017    | 4.406  | ng    | # 82      |
| 83) Chrysene                | 19.27 | 228  | 81865    | 5.516  | ng    | # 88      |
| 86) Indeno(1,2,3-cd)pyrene  | 22.64 | 276  | 45538    | 2.604  | ng    | 97        |
| 88) Benzo(b)fluoranthene    | 20.53 | 252  | 91905m   | 5.885  | ng    |           |
| 90) Benzo(a)pyrene          | 20.93 | 252  | 70543    | 4.904  | ng    | # 84      |
| 92) Benzo(g,h,i)perylene    | 23.13 | 276  | 44822    | 3.225  | ng    | # 91      |

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

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