

Data Path : Z:\HPCHEM1\BNA F\DATA\BF013118\  
 Data File : BF102666.D  
 Acq On : 31 Jan 2018 20:30  
 Operator : SJ/JU  
 Sample : J1015-26 2X  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CL-02-013018-D

Manual Integrations  
 APPROVED

Sohil  
 2/1/2018 12:14:12 PM

Quant Time: Feb 01 00:38:17 2018  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jan 31 18:38:00 2018  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) | 2/1/2018 12:14:12 PM |
|-----------------------------|-------|------|----------|--------|-------|----------|----------------------|
| 1) 1,4-Dichlorobenzene-d4   | 6.70  | 152  | 159650   | 20.00  | ng    | 0.00     |                      |
| 21) Naphthalene-d8          | 7.99  | 136  | 667581   | 20.00  | ng    | 0.00     |                      |
| 38) Acenaphthene-d10        | 9.73  | 164  | 282476   | 20.00  | ng    | 0.00     |                      |
| 63) Phenanthrene-d10        | 11.22 | 188  | 403717   | 20.00  | ng    | 0.00     |                      |
| 75) Chrysene-d12            | 13.86 | 240  | 299207   | 20.00  | ng    | 0.00     |                      |
| 86) Perylene-d12            | 15.29 | 264  | 237728   | 20.00  | ng    | 0.00     |                      |
| System Monitoring Compounds |       |      |          |        |       |          |                      |
| 5) 2-Fluorophenol           | 5.31  | 112  | 540598   | 51.34  | ng    | 0.00     |                      |
| 7) Phenol-d6                | 6.35  | 99   | 643659   | 50.71  | ng    | 0.00     |                      |
| 23) Nitrobenzene-d5         | 7.27  | 82   | 384676   | 33.11  | ng    | -0.01    |                      |
| 41) 2,4,6-Tribromophenol    | 10.53 | 330  | 107804   | 38.52  | ng    | 0.00     |                      |
| 44) 2-Fluorobiphenyl        | 9.06  | 172  | 685492   | 35.68  | ng    | 0.00     |                      |
| 78) Terphenyl-d14           | 12.80 | 244  | 442591   | 33.35  | ng    | 0.00     |                      |
| Target Compounds            |       |      |          |        |       |          |                      |
| 49) Dimethylphthalate       | 9.46  | 163  | 97773    | 4.188  | ng    |          | Qvalue # 99          |
| 70) Phenanthrene            | 11.25 | 178  | 132441   | 5.630  | ng    |          | 97                   |
| 71) Anthracene              | 11.30 | 178  | 50951    | 2.122  | ng    |          | 97                   |
| 74) Fluoranthene            | 12.43 | 202  | 327245   | 13.641 | ng    |          | 99                   |
| 77) Pyrene                  | 12.66 | 202  | 319373   | 13.806 | ng    |          | 99                   |
| 80) Benzo(a)anthracene      | 13.85 | 228  | 186144   | 10.105 | ng    |          | 97                   |
| 82) Chrysene                | 13.89 | 228  | 167026   | 8.929  | ng    |          | 97                   |
| 85) Indeno(1,2,3-cd)pyrene  | 16.69 | 276  | 61556    | 3.984  | ng    |          | 98                   |
| 87) Benzo(b)fluoranthene    | 14.89 | 252  | 198853m  | 12.906 | ng    |          |                      |
| 88) Benzo(k)fluoranthene    | 14.91 | 252  | 48293m   | 3.317  | ng    |          |                      |
| 89) Benzo(a)pyrene          | 15.23 | 252  | 126015   | 9.068  | ng    |          | # 87                 |
| 91) Benzo(a,h,i)perylene    | 17.12 | 276  | 59897    | 5.389  | ng    |          | 92                   |

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