

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102516\  
 Data File : BF090825.D  
 Acq On : 25 Oct 2016 23:22  
 Operator : UM/SJ  
 Sample : H5394-03  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SIDEWELL-NORTH

Manual Integrations  
 APPROVED

umangi  
 10/26/2016 4:46:11 PM

Quant Time: Oct 26 01:30:28 2016  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102116.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Oct 25 18:56:36 2016  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min)       |
|-----------------------------|-------|------|----------|--------|-------|----------------|
| 1) 1,4-Dichlorobenzene-d4   | 6.80  | 152  | 206137   | 20.00  | ng    | 0.00           |
| 21) Naphthalene-d8          | 8.09  | 136  | 814791   | 20.00  | ng    | 0.00           |
| 38) Acenaphthene-d10        | 9.84  | 164  | 414597   | 20.00  | ng    | -0.01          |
| 63) Phenanthrene-d10        | 11.32 | 188  | 714457   | 20.00  | ng    | 0.00           |
| 75) Chrysene-d12            | 13.96 | 240  | 476561   | 20.00  | ng    | 0.00           |
| 86) Perylene-d12            | 15.38 | 264  | 381607   | 20.00  | ng    | 0.00           |
| System Monitoring Compounds |       |      |          |        |       |                |
| 5) 2-Fluorophenol           | 5.40  | 112  | 1633770  | 127.82 | ng    | 0.01           |
| 7) Phenol-d6                | 6.44  | 99   | 1940985  | 112.11 | ng    | 0.00           |
| 23) Nitrobenzene-d5         | 7.37  | 82   | 1315841  | 88.57  | ng    | 0.00           |
| 41) 2,4,6-Tribromophenol    | 10.64 | 330  | 705279   | 214.92 | ng    | 0.00           |
| 44) 2-Fluorobiphenyl        | 9.16  | 172  | 2181868  | 81.70  | ng    | -0.01          |
| 78) Terphenyl-d14           | 12.92 | 244  | 2224397  | 106.98 | ng    | 0.00           |
| Target Compounds            |       |      |          |        |       |                |
| 49) Dimethylphthalate       | 9.56  | 163  | 669347   | 25.69  | ng    | Qvalue<br># 98 |
| 70) Phenanthrene            | 11.34 | 178  | 148258   | 4.07   | ng    | 98             |
| 74) Fluoranthene            | 12.53 | 202  | 312734   | 8.88   | ng    | 93             |
| 77) Pyrene                  | 12.76 | 202  | 242853   | 7.30   | ng    | 97             |
| 80) Benzo(a)anthracene      | 13.95 | 228  | 89547m   | 3.43   | ng    |                |
| 82) Chrysene                | 13.99 | 228  | 108608m  | 4.27   | ng    |                |
| 85) Indeno(1,2,3-cd)pyrene  | 16.74 | 276  | 52811    | 2.38   | ng    | # 93           |
| 87) Benzo(b)fluoranthene    | 14.98 | 252  | 115178m  | 4.91   | ng    |                |
| 88) Benzo(k)fluoranthene    | 15.00 | 252  | 49743m   | 2.11   | ng    |                |
| 89) Benzo(a)pyrene          | 15.32 | 252  | 72533    | 3.29   | ng    | # 86           |
| 91) Benzo(g,h,i)perylene    | 17.15 | 276  | 50513    | 2.46   | ng    | # 81           |

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

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