

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092218\  
 Data File : BF109248.D  
 Acq On : 23 Sep 2018 4:12  
 Operator : JU/SJ  
 Sample : J5017-03  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 RA-091918-N-EW-057

Manual Integrations  
 APPROVED

Sohil  
 9/24/2018 11:39:18 AM

Quant Time: Sep 24 08:26:08 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Sep 22 13:32:18 2018  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.70  | 152  | 100814   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 7.99  | 136  | 362145   | 20.00  | ng    | 0.00     |
| 38) Acenaphthene-d10        | 9.73  | 164  | 148734   | 20.00  | ng    | 0.00     |
| 63) Phenanthrene-d10        | 11.21 | 188  | 219762   | 20.00  | ng    | 0.00     |
| 75) Chrysene-d12            | 13.84 | 240  | 201516   | 20.00  | ng    | 0.00     |
| 86) Perylene-d12            | 15.24 | 264  | 171593   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.32  | 112  | 724694   | 118.40 | ng    | 0.00     |
| 7) Phenol-d6                | 6.35  | 99   | 944675   | 120.95 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.27  | 82   | 586790   | 95.65  | ng    | -0.01    |
| 41) 2,4,6-Tribromophenol    | 10.52 | 330  | 161484   | 110.14 | ng    | -0.01    |
| 44) 2-Fluorobiphenyl        | 9.06  | 172  | 1002546  | 101.66 | ng    | -0.01    |
| 78) Terphenyl-d14           | 12.80 | 244  | 680381   | 82.86  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       |          |
| 49) Dimethylphthalate       | 9.46  | 163  | 201941   | 18.667 | ng    | 98       |
| 70) Phenanthrene            | 11.23 | 178  | 125549   | 11.442 | ng    | 99       |
| 71) Anthracene              | 11.29 | 178  | 35568    | 3.196  | ng    | 96       |
| 74) Fluoranthene            | 12.42 | 202  | 119709   | 10.209 | ng    | 95       |
| 77) Pyrene                  | 12.65 | 202  | 124851   | 8.645  | ng    | 98       |
| 80) Benzo(a)anthracene      | 13.83 | 228  | 66451    | 5.475  | ng    | 96       |
| 82) Chrysene                | 13.86 | 228  | 62712m   | 5.213  | ng    |          |
| 87) Benzo(b)fluoranthene    | 14.84 | 252  | 50965m   | 5.405  | ng    |          |
| 89) Benzo(a)pyrene          | 15.18 | 252  | 35038    | 4.124  | ng    | # 87     |
| 91) Benzo(g,h,i)perylene    | 16.97 | 276  | 14758    | 2.154  | ng    | 93       |

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

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