

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080917\  
 Data File : BF097544.D  
 Acq On : 10 Aug 2017 4:30  
 Operator : SJ/JU  
 Sample : I4646-05  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 E2-L10-BASE

Manual Integrations  
 APPROVED

Sohil  
 8/10/2017 6:08:02 PM

Quant Time: Aug 10 08:00:28 2017  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 07 16:02:51 2017  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min)   |
|-----------------------------|-------|------|----------|--------|-------|------------|
| 1) 1,4-Dichlorobenzene-d4   | 6.40  | 152  | 100122   | 20.00  | ng    | -0.01      |
| 21) Naphthalene-d8          | 7.68  | 136  | 449202   | 20.00  | ng    | -0.02      |
| 38) Acenaphthene-d10        | 9.42  | 164  | 181969   | 20.00  | ng    | -0.02      |
| 63) Phenanthrene-d10        | 10.89 | 188  | 272414   | 20.00  | ng    | -0.02      |
| 75) Chrysene-d12            | 13.52 | 240  | 192544   | 20.00  | ng    | -0.02      |
| 86) Perylene-d12            | 14.86 | 264  | 180935   | 20.00  | ng    | -0.01      |
| System Monitoring Compounds |       |      |          |        |       |            |
| 5) 2-Fluorophenol           | 4.99  | 112  | 529507   | 79.73  | ng    | -0.02      |
| 7) Phenol-d6                | 6.07  | 99   | 654779   | 86.36  | ng    | -0.02      |
| 23) Nitrobenzene-d5         | 6.97  | 82   | 408071   | 53.29  | ng    | -0.02      |
| 41) 2,4,6-Tribromophenol    | 10.21 | 330  | 106151   | 73.83  | ng    | -0.02      |
| 44) 2-Fluorobiphenyl        | 8.76  | 172  | 673179   | 62.78  | ng    | -0.02      |
| 78) Terphenyl-d14           | 12.47 | 244  | 405389   | 42.93  | ng    | -0.02      |
| Target Compounds            |       |      |          |        |       |            |
| 49) Dimethylphthalate       | 9.16  | 163  | 133708   | 9.676  | ng    | Qvalue 100 |
| 70) Phenanthrene            | 10.92 | 178  | 157620   | 10.799 | ng    | 97         |
| 71) Anthracene              | 10.97 | 178  | 30323m   | 2.028  | ng    |            |
| 74) Fluoranthene            | 12.10 | 202  | 167957   | 11.746 | ng    | 98         |
| 77) Pyrene                  | 12.32 | 202  | 140443   | 8.910  | ng    | 99         |
| 80) Benzo(a)anthracene      | 13.51 | 228  | 57827    | 4.642  | ng    | 98         |
| 82) Chrysene                | 13.54 | 228  | 50788    | 4.175  | ng    | 98         |
| 85) Indeno(1,2,3-cd)pyrene  | 16.05 | 276  | 23769    | 2.279  | ng    | 98         |
| 87) Benzo(b)fluoranthene    | 14.52 | 252  | 60478m   | 5.560  | ng    |            |
| 88) Benzo(k)fluoranthene    | 14.53 | 252  | 23939m   | 2.310  | ng    |            |
| 89) Benzo(a)pyrene          | 14.81 | 252  | 45008    | 4.521  | ng    | # 95       |
| 91) Benzo(a,h,i)perylene    | 16.40 | 276  | 20267    | 2.368  | ng    | # 89       |

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

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