

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012017\  
 Data File : BF092373.D  
 Acq On : 20 Jan 2017 15:21  
 Operator : UM/SJ  
 Sample : I1304-04  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 1634-NOTTINGHAM-COMP

Manual Integrations  
 APPROVED

mohammad  
 1/23/2017 2:44:47 PM

Quant Time: Jan 20 23:10:20 2017  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jan 20 22:22:14 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min)       |
|--------------------------------|-------|------|----------|--------|-------|----------------|
| 1) 1,4-Dichlorobenzene-d4      | 6.52  | 152  | 150615   | 20.00  | ng    | 0.00           |
| 21) Naphthalene-d8             | 7.81  | 136  | 578221   | 20.00  | ng    | 0.00           |
| 38) Acenaphthene-d10           | 9.56  | 164  | 438504   | 20.00  | ng    | 0.00           |
| 63) Phenanthrene-d10           | 11.03 | 188  | 667979   | 20.00  | ng    | 0.00           |
| 75) Chrysene-d12               | 13.66 | 240  | 412790   | 20.00  | ng    | 0.00           |
| 86) Perylene-d12               | 15.01 | 264  | 364469   | 20.00  | ng    | 0.00           |
| System Monitoring Compounds    |       |      |          |        |       |                |
| 5) 2-Fluorophenol              | 5.11  | 112  | 1377817  | 152.75 | ng    | 0.02           |
| 7) Phenol-d6                   | 6.20  | 99   | 1843668  | 167.41 | ng    | 0.01           |
| 23) Nitrobenzene-d5            | 7.10  | 82   | 1155943  | 111.16 | ng    | 0.00           |
| 41) 2,4,6-Tribromophenol       | 10.35 | 330  | 425986   | 117.39 | ng    | 0.00           |
| 44) 2-Fluorobiphenyl           | 8.90  | 172  | 2188593  | 105.35 | ng    | 0.00           |
| 78) Terphenyl-d14              | 12.62 | 244  | 1611073  | 94.66  | ng    | 0.00           |
| Target Compounds               |       |      |          |        |       |                |
| 10) Phenol                     | 6.20  | 94   | 61184    | 4.88   | ng    | Qvalue<br># 14 |
| 49) Dimethylphthalate          | 9.30  | 163  | 365835   | 13.04  | ng    | 99             |
| 70) Phenanthrene               | 11.06 | 178  | 91128    | 2.52   | ng    | 97             |
| 74) Fluoranthene               | 12.23 | 202  | 200496   | 2.72   | ng    | 97             |
| 77) Pyrene                     | 12.46 | 202  | 188215   | 6.21   | ng    | 98             |
| 80) Benzo(a)anthracene         | 13.65 | 228  | 71569m   | 3.07   | ng    |                |
| 82) Chrysene                   | 13.69 | 228  | 82729m   | 3.54   | ng    |                |
| 83) Bis(2-ethylhexyl)phthalate | 13.66 | 149  | 154657   | 9.53   | ng    | 99             |
| 85) Indeno(1,2,3-cd)pyrene     | 16.23 | 276  | 42917    | 2.12   | ng    | 94             |
| 87) Benzo(b)fluoranthene       | 14.66 | 252  | 88919m   | 3.92   | ng    |                |
| 89) Benzo(a)pyrene             | 14.97 | 252  | 62950    | 3.13   | ng    | 98             |
| 91) Benzo(a,h,i)perylene       | 16.59 | 276  | 42660    | 2.48   | ng    | # 91           |

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

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