

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062317\  
 Data File : BF096172.D  
 Acq On : 23 Jun 2017 22:03  
 Operator : SJ/MA  
 Sample : I3743-05  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 FNDFL-VCTY-061517-COMP

Manual Integrations  
 APPROVED

mohammad  
 6/28/2017 1:11:38 PM

Quant Time: Jun 24 00:57:01 2017  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 20 16:05:47 2017  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.26  | 152  | 71028    | 20.00 | ng    | -0.04    |
| 21) Naphthalene-d8          | 9.29  | 136  | 294484   | 20.00 | ng    | -0.04    |
| 38) Acenaphthene-d10        | 12.11 | 164  | 127768   | 20.00 | ng    | -0.04    |
| 63) Phenanthrene-d10        | 14.50 | 188  | 215661   | 20.00 | ng    | -0.04    |
| 75) Chrysene-d12            | 18.24 | 240  | 149081   | 20.00 | ng    | -0.03    |
| 86) Perylene-d12            | 19.92 | 264  | 144108   | 20.00 | ng    | -0.02    |
| System Monitoring Compounds |       |      |          |       |       |          |
| 5) 2-Fluorophenol           | 5.15  | 112  | 418630   | 93.92 | ng    | -0.04    |
| 7) Phenol-d6                | 6.84  | 99   | 523051   | 98.02 | ng    | -0.04    |
| 23) Nitrobenzene-d5         | 8.18  | 82   | 305485   | 64.42 | ng    | -0.04    |
| 41) 2,4,6-Tribromophenol    | 13.41 | 330  | 101676   | 89.94 | ng    | -0.04    |
| 44) 2-Fluorobiphenyl        | 11.07 | 172  | 606349   | 66.04 | ng    | -0.04    |
| 78) Terphenyl-d14           | 16.89 | 244  | 399440   | 66.49 | ng    | -0.04    |
| Target Compounds            |       |      |          |       |       | Qvalue   |
| 49) Dimethylphthalate       | 11.77 | 163  | 80829    | 8.33  | ng    | 99       |
| 70) Phenanthrene            | 14.55 | 178  | 41033m   | 3.30  | ng    |          |
| 74) Fluoranthene            | 16.34 | 202  | 67333    | 5.47  | ng    | 98       |
| 77) Pyrene                  | 16.64 | 202  | 61330    | 5.51  | ng    | 97       |
| 80) Benzo(a)anthracene      | 18.23 | 228  | 33668    | 3.88  | ng    | 97       |
| 82) Chrysene                | 18.28 | 228  | 31885    | 3.68  | ng    | 91       |
| 85) Indeno(1,2,3-cd)pyrene  | 21.10 | 276  | 22433    | 3.03  | ng    | 93       |
| 87) Benzo(b)fluoranthene    | 19.52 | 252  | 52280    | 5.79  | ng    | # 96     |
| 89) Benzo(a)pyrene          | 19.87 | 252  | 35105    | 4.23  | ng    | 95       |
| 91) Benzo(g,h,i)perylene    | 21.42 | 276  | 19779    | 2.76  | ng    | 93       |

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

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