

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100215\  
 Data File : BF082040.D  
 Acq On : 4 Oct 2015 00:24  
 Operator : UM/IZ  
 Sample : G3891-09  
 Misc :  
 ALS Vial : 60 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SS-9(0-2)

Manual Integrations  
 APPROVED

MMDADODA  
 10/5/2015 6:27:34 PM

Quant Time: Oct 05 04:52:13 2015  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Sep 28 18:10:42 2015  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.89  | 152  | 76740    | 20.00  | ng    | -0.05    |
| 21) Naphthalene-d8          | 8.17  | 136  | 317234   | 20.00  | ng    | -0.05    |
| 38) Acenaphthene-d10        | 9.93  | 164  | 155896   | 20.00  | ng    | -0.05    |
| 63) Phenanthrene-d10        | 11.43 | 188  | 290477   | 20.00  | ng    | -0.03    |
| 75) Chrysene-d12            | 14.07 | 240  | 238645   | 20.00  | ng    | -0.05    |
| 86) Perylene-d12            | 15.53 | 264  | 221835   | 20.00  | ng    | -0.05    |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.49  | 112  | 480363   | 113.63 | ng    | -0.03    |
| 7) Phenol-d6                | 6.51  | 99   | 641194   | 120.54 | ng    | -0.05    |
| 23) Nitrobenzene-d5         | 7.45  | 82   | 377245   | 79.27  | ng    | -0.05    |
| 41) 2,4,6-Tribromophenol    | 10.73 | 330  | 163027   | 103.26 | ng    | -0.03    |
| 44) 2-Fluorobiphenyl        | 9.26  | 172  | 783855   | 83.93  | ng    | -0.05    |
| 78) Terphenyl-d14           | 13.02 | 244  | 687850   | 96.75  | ng    | -0.03    |
| Target Compounds            |       |      |          |        |       | Qvalue   |
| 37) 2-Methylnaphthalene     | 8.88  | 142  | 21607    | 2.25   | ng    | # 88     |
| 49) Dimethylphthalate       | 9.65  | 163  | 165088   | 15.34  | ng    | # 97     |
| 70) Phenanthrene            | 11.45 | 178  | 53999    | 2.28   | ng    | 97       |
| 74) Fluoranthene            | 12.64 | 202  | 82129    | 5.28   | ng    | 88       |
| 77) Pyrene                  | 12.87 | 202  | 73272    | 5.14   | ng    | 97       |
| 80) Benzo(a)anthracene      | 14.06 | 228  | 44049    | 3.43   | ng    | # 86     |
| 82) Chrysene                | 14.09 | 228  | 41398m   | 3.33   | ng    |          |
| 85) Indeno(1,2,3-cd)pyrene  | 16.98 | 276  | 24392m   | 2.43   | ng    |          |
| 87) Benzo(b)fluoranthene    | 15.11 | 252  | 59336m   | 4.68   | ng    |          |
| 88) Benzo(k)fluoranthene    | 15.12 | 252  | 31415m   | 2.71   | ng    |          |
| 89) Benzo(a)pyrene          | 15.48 | 252  | 39632    | 3.61   | ng    | # 87     |
| 91) Benzo(a,h,i)perylene    | 17.42 | 276  | 27061    | 2.86   | ng    | # 80     |

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

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