

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\  
 Data File : BF116695.D  
 Acq On : 31 Aug 2019 13:23  
 Operator : HP/JU  
 Sample : K4643-19  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 T11-D1-D4-(0-4)

Manual Integrations  
 APPROVED

mohammad  
 9/5/2019 9:23:22 AM

Quant Time: Aug 31 17:43:52 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Aug 30 20:02:30 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4      | 6.85  | 152  | 108367   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8             | 8.13  | 136  | 425696   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10           | 9.88  | 164  | 220117   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10           | 11.36 | 188  | 358940   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12               | 13.99 | 240  | 222281   | 20.00  | ng    | 0.00     |
| 87) Perylene-d12               | 15.44 | 264  | 255527   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds    |       |      |          |        |       |          |
| 5) 2-Fluorophenol              | 5.46  | 112  | 556457   | 83.19  | ng    | 0.00     |
| 7) Phenol-d6                   | 6.47  | 99   | 764502   | 87.04  | ng    | 0.00     |
| 23) Nitrobenzene-d5            | 7.40  | 82   | 475255   | 63.73  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol       | 10.66 | 330  | 137902   | 93.71  | ng    | -0.01    |
| 45) 2-Fluorobiphenyl           | 9.20  | 172  | 791521   | 60.66  | ng    | 0.00     |
| 79) Terphenyl-d14              | 12.94 | 244  | 708957   | 59.00  | ng    | 0.00     |
| Target Compounds               |       |      |          |        |       |          |
| 50) Dimethylphthalate          | 9.59  | 163  | 93473    | 6.151  | ng    | 99       |
| 71) Phenanthrene               | 11.39 | 178  | 136353   | 6.824  | ng    | 99       |
| 75) Fluoranthene               | 12.57 | 202  | 286742   | 14.603 | ng    | 96       |
| 78) Pyrene                     | 12.80 | 202  | 234299   | 11.858 | ng    | 97       |
| 81) Benzo(a)anthracene         | 13.98 | 228  | 96464    | 6.857  | ng    | 96       |
| 83) Chrysene                   | 14.02 | 228  | 131659m  | 9.084  | ng    |          |
| 84) Bis(2-ethylhexyl)phthalate | 13.98 | 149  | 77456    | 6.530  | ng    | # 95     |
| 86) Indeno(1,2,3-cd)pyrene     | 16.88 | 276  | 84278    | 7.239  | ng    | 99       |
| 88) Benzo(b)fluoranthene       | 15.02 | 252  | 166165m  | 10.756 | ng    |          |
| 89) Benzo(k)fluoranthene       | 15.05 | 252  | 59558m   | 4.228  | ng    |          |
| 90) Benzo(a)pyrene             | 15.38 | 252  | 99876    | 7.431  | ng    | 97       |
| 92) Benzo(a,h,i)perylene       | 17.32 | 276  | 81824    | 6.820  | ng    | # 92     |

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

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