

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070219\  
 Data File : BF115361.D  
 Acq On : 2 Jul 2019 16:00  
 Operator : HP/JU  
 Sample : K3606-02DL 5X  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 HJDC-COMP-1DL

Manual Integrations  
 APPROVED

mohammad  
 7/3/2019 10:51:36 PM

Quant Time: Jul 02 16:33:59 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jul 02 16:24:08 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.61  | 152  | 192749   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 7.89  | 136  | 779020   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.64  | 164  | 388667   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.11 | 188  | 763817   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.74 | 240  | 446114   | 20.00  | ng    | -0.01    |
| 87) Perylene-d12            | 15.12 | 264  | 503963   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.20  | 112  | 172377   | 13.80  | ng    | 0.00     |
| 7) Phenol-d6                | 6.24  | 99   | 232370   | 15.33  | ng    | -0.02    |
| 23) Nitrobenzene-d5         | 7.16  | 82   | 128771   | 10.17  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol    | 10.42 | 330  | 59055    | 14.41  | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 8.97  | 172  | 295393   | 12.91  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.70 | 244  | 313039   | 14.92  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 11.14 | 178  | 791757   | 19.252 | ng    | 97       |
| 72) Anthracene              | 11.19 | 178  | 270994   | 6.528  | ng    | 97       |
| 75) Fluoranthene            | 12.32 | 202  | 1437317  | 35.029 | ng    | 99       |
| 78) Pyrene                  | 12.55 | 202  | 1399561  | 40.822 | ng    | 98       |
| 81) Benzo(a)anthracene      | 13.74 | 228  | 582876   | 18.209 | ng    | 99       |
| 83) Chrysene                | 13.77 | 228  | 505238   | 17.231 | ng    | 97       |
| 86) Indeno(1,2,3-cd)pyrene  | 16.39 | 276  | 313840   | 9.045  | ng    | 95       |
| 88) Benzo(b)fluoranthene    | 14.74 | 252  | 665004m  | 21.148 | ng    |          |
| 89) Benzo(k)fluoranthene    | 14.77 | 252  | 151676m  | 5.379  | ng    |          |
| 90) Benzo(a)pyrene          | 15.07 | 252  | 494211   | 17.437 | ng    | 98       |
| 91) Dibenzo(a,h)anthracene  | 16.41 | 278  | 77454m   | 2.927  | ng    |          |
| 92) Benzo(a,h,i)perylene    | 16.78 | 276  | 283939   | 10.603 | ng    | 99       |

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

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