

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022620\  
 Data File : BF119095.D  
 Acq On : 26 Feb 2020 15:24  
 Operator : CG/JU  
 Sample : L1678-01  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 VNJ-241

Manual Integrations  
 APPROVED

mohammad  
 2/27/2020 4:58:30 PM

Quant Time: Feb 26 16:39:41 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Feb 26 16:03:05 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.75  | 152  | 158254   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.03  | 136  | 601142   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.77  | 164  | 312505   | 20.00  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.26 | 188  | 508965   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.89 | 240  | 317035   | 20.00  | ng    | -0.01    |
| 87) Perylene-d12            | 15.32 | 264  | 374546   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.37  | 112  | 876281   | 92.54  | ng    | 0.00     |
| 7) Phenol-d6                | 6.39  | 99   | 1064314  | 92.41  | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.31  | 82   | 710511   | 62.41  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.56 | 330  | 337317   | 82.51  | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.10  | 172  | 1325746  | 62.50  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.84 | 244  | 1151008  | 62.51  | ng    | -0.01    |
| Target Compounds            |       |      |          |        |       | Qvalue   |
| 50) Dimethylphthalate       | 9.49  | 163  | 93555    | 3.915  | ng    | 99       |
| 71) Phenanthrene            | 11.28 | 178  | 367729   | 13.248 | ng    | 98       |
| 72) Anthracene              | 11.33 | 178  | 102530   | 3.697  | ng    | 97       |
| 75) Fluoranthene            | 12.47 | 202  | 649697   | 22.052 | ng    | 98       |
| 78) Pyrene                  | 12.70 | 202  | 521455   | 20.483 | ng    | 98       |
| 81) Benzo(a)anthracene      | 13.89 | 228  | 207693   | 9.615  | ng    | 99       |
| 83) Chrysene                | 13.92 | 228  | 197725   | 9.186  | ng    | 97       |
| 86) Indeno(1,2,3-cd)pyrene  | 16.72 | 276  | 113684   | 5.455  | ng    | 94       |
| 88) Benzo(b)fluoranthene    | 14.92 | 252  | 240951m  | 9.760  | ng    |          |
| 89) Benzo(k)fluoranthene    | 14.94 | 252  | 72267m   | 3.159  | ng    |          |
| 90) Benzo(a)pyrene          | 15.26 | 252  | 162978   | 7.314  | ng    | # 97     |
| 92) Benzo(a,h,i)perylene    | 17.14 | 276  | 111520   | 5.757  | ng    | 94       |

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

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