

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF060120\  
 Data File : BF120466.D  
 Acq On : 1 Jun 2020 13:04  
 Operator : JU/CG  
 Sample : L2792-01  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 NM32-COMP-2020-0-6

Manual Integrations  
 APPROVED

mohammad  
 6/2/2020 3:24:32 PM

Quant Time: Jun 01 13:48:43 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF052820.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri May 29 12:39:43 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.84  | 152  | 242982   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.12  | 136  | 863465   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.87  | 164  | 397111   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.36 | 188  | 638262   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.00 | 240  | 494543   | 20.00  | ng    | 0.00     |
| 86) Perylene-d12            | 15.45 | 264  | 565234   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.46  | 112  | 1267266  | 99.39  | ng    | 0.00     |
| 7) Phenol-d6                | 6.47  | 99   | 1613283  | 102.66 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.40  | 82   | 1052410  | 79.73  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.66 | 330  | 385989   | 104.91 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.20  | 172  | 1854295  | 79.23  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.94 | 244  | 1622308  | 73.00  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       | Qvalue   |
| 50) Dimethylphthalate       | 9.59  | 163  | 249034   | 10.03  | ng    | 99       |
| 71) Phenanthrene            | 11.38 | 178  | 188796   | 6.71   | ng    | 97       |
| 75) Fluoranthene            | 12.57 | 202  | 406817   | 13.44  | ng    | 97       |
| 78) Pyrene                  | 12.80 | 202  | 384189   | 11.17  | ng    | 98       |
| 81) Benzo(a)anthracene      | 13.99 | 228  | 219530   | 8.28   | ng    | 97       |
| 83) Chrysene                | 14.02 | 228  | 211105   | 7.60   | ng    | 97       |
| 87) Indeno(1,2,3-cd)pyrene  | 16.89 | 276  | 152969m  | 4.95   | ng    |          |
| 88) Benzo(b)fluoranthene    | 15.03 | 252  | 326003m  | 10.38  | ng    |          |
| 89) Benzo(k)fluoranthene    | 15.05 | 252  | 96763m   | 3.30   | ng    |          |
| 90) Benzo(a)pyrene          | 15.39 | 252  | 234209   | 8.53   | ng    | 99       |
| 92) Benzo(g,h,i)perylene    | 17.32 | 276  | 137652   | 5.54   | ng    | 96       |

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

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