

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101922\  
 Data File : BF130723.D  
 Acq On : 19 Oct 2022 21:31  
 Operator : CG\JU  
 Sample : N5175-01  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CONCRETE-1

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 10/20/2022  
 Supervised By : mohammad ahmed 10/20/2022

Quant Time: Oct 20 02:47:40 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Oct 18 03:15:54 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.569  | 152  | 76990    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 7.851  | 136  | 353069   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.598  | 164  | 157210   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.080 | 188  | 238872   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.710 | 240  | 170451   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 15.086 | 264  | 165960   | 20.000  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.175  | 112  | 399339   | 93.496  | ng    | 0.00     |
| 7) Phenol-d6                | 6.228  | 99   | 671483   | 125.042 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.139  | 82   | 378006   | 58.446  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.392 | 330  | 168512   | 109.545 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 8.928  | 172  | 886616   | 85.427  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.663 | 244  | 903705   | 87.996  | ng    | -0.01    |
|                             |        |      |          |         |       |          |
| Target Compounds            |        |      |          |         |       | Qvalue   |
| 52) Acenaphthene            | 9.628  | 154  | 29623    | 3.504   | ng    | 98       |
| 58) Fluorene                | 10.145 | 166  | 32753    | 3.094   | ng    | 99       |
| 71) Phenanthrene            | 11.104 | 178  | 410392   | 31.010  | ng    | 100      |
| 72) Anthracene              | 11.157 | 178  | 101794   | 7.876   | ng    | 98       |
| 75) Fluoranthene            | 12.286 | 202  | 457181   | 35.943  | ng    | 97       |
| 78) Pyrene                  | 12.516 | 202  | 521849   | 35.488  | ng    | 99       |
| 81) Benzo(a)anthracene      | 13.698 | 228  | 236364   | 20.580  | ng    | 98       |
| 83) Chrysene                | 13.739 | 228  | 233205   | 21.416  | ng    | 98       |
| 87) Indeno(1,2,3-cd)pyrene  | 16.380 | 276  | 103258   | 8.313   | ng    | 97       |
| 88) Benzo(b)fluoranthene    | 14.710 | 252  | 258472m  | 24.595  | ng    |          |
| 89) Benzo(k)fluoranthene    | 14.733 | 252  | 80878m   | 7.558   | ng    |          |
| 90) Benzo(a)pyrene          | 15.033 | 252  | 199137   | 22.768  | ng    | 97       |
| 91) Dibenzo(a,h)anthracene  | 16.380 | 278  | 29092    | 2.857   | ng    | # 86     |
| 92) Benzo(g,h,i)perylene    | 16.768 | 276  | 80397    | 8.478   | ng    | 99       |
| -----                       |        |      |          |         |       |          |

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

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