

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF092023\  
 Data File : BF135364.D  
 Acq On : 20 Sep 2023 18:43  
 Operator : CG\JU  
 Sample : 04397-04  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 B103-04

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 09/21/2023  
 Supervised By :mohammad ahmed 09/21/2023

Quant Time: Sep 21 01:52:47 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Sep 12 00:19:07 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.757  | 152  | 129628   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.034  | 136  | 475315   | 20.000 | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.792  | 164  | 206218   | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.286 | 188  | 319985   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 13.939 | 240  | 250380   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 15.410 | 264  | 264513   | 20.000 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.393  | 112  | 660443   | 82.866 | ng    | 0.00     |
| 7) Phenol-d6                | 6.398  | 99   | 812777   | 80.200 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.322  | 82   | 485007   | 55.437 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.586 | 330  | 123637   | 60.331 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.116  | 172  | 710550   | 51.985 | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.880 | 244  | 506945   | 31.387 | ng    | -0.01    |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 11.310 | 178  | 205976   | 13.610 | ng    | 98       |
| 72) Anthracene              | 11.363 | 178  | 41216    | 2.650  | ng    | 97       |
| 75) Fluoranthene            | 12.504 | 202  | 294111   | 18.651 | ng    | 99       |
| 78) Pyrene                  | 12.733 | 202  | 271096   | 11.372 | ng    | 99       |
| 81) Benzo(a)anthracene      | 13.933 | 228  | 142205   | 8.796  | ng    | 98       |
| 83) Chrysene                | 13.968 | 228  | 136591   | 8.497  | ng    | 97       |
| 87) Indeno(1,2,3-cd)pyrene  | 16.886 | 276  | 44341    | 2.557  | ng    | 97       |
| 88) Benzo(b)fluoranthene    | 14.980 | 252  | 157509m  | 11.000 | ng    |          |
| 89) Benzo(k)fluoranthene    | 15.004 | 252  | 56702m   | 4.009  | ng    |          |
| 90) Benzo(a)pyrene          | 15.345 | 252  | 104616   | 7.661  | ng    | # 97     |
| 92) Benzo(g,h,i)perylene    | 17.333 | 276  | 41740    | 2.816  | ng    | # 93     |
| -----                       |        |      |          |        |       |          |

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

## Manual Integrations

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