

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF081623\  
 Data File : BF134815.D  
 Acq On : 16 Aug 2023 18:36  
 Operator : CG\JU  
 Sample : 03994-01DL2 500X  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 NEVINS-SB06-Z22-23DL2

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 08/17/2023  
 Supervised By :mohammad ahmed 08/17/2023

Quant Time: Aug 17 01:36:15 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF080123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Aug 04 09:29:15 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.798  | 152  | 143775   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.080  | 136  | 561687   | 20.000 | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.833  | 164  | 242134   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.321 | 188  | 351957   | 20.000 | ng    | # 0.00   |
| 76) Chrysene-d12            | 13.974 | 240  | 244884   | 20.000 | ng    | # 0.00   |
| 86) Perylene-d12            | 15.445 | 264  | 179532   | 20.000 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 0.000  | 112  | 0d       | 0.000  | ng    |          |
| 7) Phenol-d6                | 6.433  | 99   | 720      | 0.060  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 0.000  | 82   | 0d       | 0.000  | ng    |          |
| 42) 2,4,6-Tribromophenol    | 0.000  | 330  | 0        | 0.000  | ng    |          |
| 45) 2-Fluorobiphenyl        | 9.151  | 172  | 1417     | 0.097  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.915 | 244  | 1682     | 0.121  | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 31) Naphthalene             | 8.104  | 128  | 2079741  | 73.673 | ng    | 98       |
| 37) 2-Methylnaphthalene     | 8.792  | 142  | 650265   | 34.752 | ng    | 98       |
| 38) 1-Methylnaphthalene     | 8.892  | 142  | 401933   | 23.100 | ng    | 97       |
| 46) 1,1'-Biphenyl           | 9.257  | 154  | 97884    | 5.199  | ng    | 96       |
| 52) Acenaphthene            | 9.868  | 154  | 282887   | 19.683 | ng    | 98       |
| 58) Fluorene                | 10.380 | 166  | 124387   | 8.347  | ng    | 95       |
| 71) Phenanthrene            | 11.345 | 178  | 457556   | 24.475 | ng    | 98       |
| 72) Anthracene              | 11.398 | 178  | 128936   | 6.752  | ng    | 98       |
| 75) Fluoranthene            | 12.539 | 202  | 170777   | 8.648  | ng    | 97       |
| 78) Pyrene                  | 12.768 | 202  | 269152   | 12.702 | ng    | 99       |
| 81) Benzo(a)anthracene      | 13.962 | 228  | 74685m   | 4.265  | ng    |          |
| 83) Chrysene                | 13.998 | 228  | 60533    | 3.548  | ng    | 98       |
| 88) Benzo(b)fluoranthene    | 15.015 | 252  | 31181m   | 2.848  | ng    |          |
| 90) Benzo(a)pyrene          | 15.380 | 252  | 38994    | 3.820  | ng    | # 93     |
| -----                       |        |      |          |        |       |          |

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

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