

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF010722\  
 Data File : BF126557.D  
 Acq On : 7 Jan 2022 22:05  
 Operator : JU/CG  
 Sample : N1068-01 5X  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TR-05-010622

Quant Time: Jan 08 04:41:10 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122921.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jan 05 01:57:07 2022  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

Reviewed By :Christian Giraldo 01/10/2022  
 Supervised By :Jagrut Upadhyay 01/10/2022

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.781  | 152  | 77295    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.063  | 136  | 290658   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.822  | 164  | 125071   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.310 | 188  | 210425   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 13.957 | 240  | 141127   | 20.000 | ng    | # 0.00   |
| 86) Perylene-d12            | 15.416 | 264  | 103540   | 20.000 | ng    | 0.03     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.387  | 112  | 95393    | 17.723 | ng    | 0.00     |
| 7) Phenol-d6                | 6.410  | 99   | 111908   | 16.052 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.340  | 82   | 64003    | 11.306 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.610 | 330  | 17931    | 18.553 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.139  | 172  | 107569   | 15.135 | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.898 | 244  | 104193   | 14.345 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 11.333 | 178  | 22890    | 2.128  | ng    | 96       |
| 75) Fluoranthene            | 12.522 | 202  | 41297    | 4.206  | ng    | 100      |
| 78) Pyrene                  | 12.751 | 202  | 39879    | 3.344  | ng    | 98       |
| 81) Benzo(a)anthracene      | 13.945 | 228  | 18993    | 2.340  | ng    | # 84     |
| 83) Chrysene                | 13.980 | 228  | 21437    | 2.575  | ng    | # 92     |
| 88) Benzo(b)fluoranthene    | 14.998 | 252  | 21840m   | 3.541  | ng    |          |
| 90) Benzo(a)pyrene          | 15.351 | 252  | 14997    | 2.900  | ng    | # 50     |
| -----                       |        |      |          |        |       |          |

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

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