

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF081922\  
 Data File : BF129922.D  
 Acq On : 19 Aug 2022 22:03  
 Operator : CG\JU  
 Sample : N4093-04  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BLDG-4-BASEMENT-DRUM

Quant Time: Aug 20 07:54:09 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF081422.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Aug 20 07:47:33 2022  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

Reviewed By :Christian Giraldo 08/22/2022  
 Supervised By :Jagrut Upadhyay 08/22/2022

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.786  | 152  | 140375   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.069  | 136  | 573383   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.822  | 164  | 316560   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.310 | 188  | 553136   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.957 | 240  | 411362   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.415 | 264  | 307032   | 20.000  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.439  | 112  | 1121752  | 132.309 | ng    | 0.02     |
| 7) Phenol-d6                | 6.445  | 99   | 1302989  | 122.725 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.357  | 82   | 971455   | 90.696  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.621 | 330  | 466455   | 146.815 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.145  | 172  | 1777342  | 85.274  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.904 | 244  | 2158600  | 87.314  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| Target Compounds            |        |      |          |         |       | Qvalue   |
| 22) Acetophenone            | 7.198  | 105  | 38659    | 2.767   | ng    | # 93     |
| 32) Benzoic acid            | 7.869  | 122  | 5826m    | 4.368   | ng    |          |
| -----                       |        |      |          |         |       |          |

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

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