

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF030322\  
 Data File : BF127193.D  
 Acq On : 04 Mar 2022 01:36  
 Operator : CG\JU  
 Sample : N1708-17DL 10X  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PPSP-B10-17.5-18.0DL

Manual Integrations  
 APPROVED

Reviewed By :Christian Giraldo 03/04/2022  
 Supervised By :Jagrut Upadhyay 03/04/2022

Quant Time: Mar 04 02:16:12 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF021622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 01 23:57:09 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.893  | 152  | 121787   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.175  | 136  | 459564   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.939  | 164  | 218621   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.428 | 188  | 315065   | 20.000 | ng    | # 0.00   |
| 76) Chrysene-d12            | 14.069 | 240  | 218130   | 20.000 | ng    | # 0.00   |
| 86) Perylene-d12            | 15.563 | 264  | 192681   | 20.000 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.504  | 112  | 113432   | 14.395 | ng    | 0.00     |
| 7) Phenol-d6                | 6.516  | 99   | 150239   | 14.130 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.451  | 82   | 108076   | 10.615 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.728 | 330  | 29489    | 11.715 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.251  | 172  | 161729   | 10.891 | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.004 | 244  | 132219   | 8.911  | ng    | -0.01    |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 31) Naphthalene             | 8.198  | 128  | 320784   | 13.372 | ng    | 96       |
| 37) 2-Methylnaphthalene     | 8.892  | 142  | 956141   | 61.422 | ng    | 97       |
| 38) 1-Methylnaphthalene     | 8.992  | 142  | 554243   | 37.181 | ng    | 99       |
| 46) 1,1'-Biphenyl           | 9.351  | 154  | 159714   | 9.221  | ng    | 98       |
| 52) Acenaphthene            | 9.969  | 154  | 44232m   | 3.260  | ng    |          |
| 58) Fluorene                | 10.486 | 166  | 102821   | 7.075  | ng    | # 89     |
| 71) Phenanthrene            | 11.451 | 178  | 253719   | 14.387 | ng    | 96       |
| -----                       |        |      |          |        |       |          |

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

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