

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF062123\  
 Data File : BF133885.D  
 Acq On : 21 Jun 2023 22:47  
 Operator : CG\JU  
 Sample : 03289-19 2X  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-07-0-2.0

Quant Time: Jun 22 06:02:06 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF060923.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 20 13:43:46 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.857  | 152  | 94508    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.140  | 136  | 307287   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.898  | 164  | 119071   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.392 | 188  | 227133   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 14.045 | 240  | 171572   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 15.557 | 264  | 120393   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.487  | 112  | 168613   | 31.060 | ng    | 0.00     |
| 7) Phenol-d6                | 6.487  | 99   | 192402   | 27.228 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.416  | 82   | 120046   | 21.278 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.686 | 330  | 40214    | 26.681 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.216  | 172  | 202949   | 24.229 | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.986 | 244  | 226480   | 20.425 | ng    | 0.00     |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 11.416 | 178  | 29640    | 2.572  | ng    | 96       |
| 75) Fluoranthene            | 12.610 | 202  | 91868    | 7.507  | ng    | 98       |
| 78) Pyrene                  | 12.839 | 202  | 89765    | 5.616  | ng    | 98       |
| 81) Benzo(a)anthracene      | 14.039 | 228  | 41156    | 3.468  | ng    | 97       |
| 83) Chrysene                | 14.074 | 228  | 39900    | 3.555  | ng    | 95       |
| 88) Benzo(b)fluoranthene    | 15.110 | 252  | 37208m   | 5.088  | ng    |          |
| 90) Benzo(a)pyrene          | 15.492 | 252  | 25068    | 3.693  | ng    | 97       |
| 92) Benzo(g,h,i)perylene    | 17.574 | 276  | 18219    | 2.680  | ng    | # 89     |
| -----                       |        |      |          |        |       |          |

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

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