

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060722\  
 Data File : BF128713.D  
 Acq On : 07 Jun 2022 13:18  
 Operator : CG\JU  
 Sample : N3189-04RE  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 20220428-AMENDED-COMPOSITE-HRE

Quant Time: Jun 08 00:25:57 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF060222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jun 02 17:07:23 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.822  | 152  | 81158    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.104  | 136  | 313869   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.863  | 164  | 185663   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.351 | 188  | 343535   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 13.998 | 240  | 243945   | 20.000 | ng    | # 0.00   |
| 86) Perylene-d12            | 15.462 | 264  | 198475   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.487  | 112  | 78987    | 15.502 | ng    | 0.01     |
| 7) Phenol-d6                | 6.487  | 99   | 436827   | 70.234 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.386  | 82   | 496461   | 73.644 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.657 | 330  | 20830    | 9.495  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.186  | 172  | 891507   | 67.263 | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.945 | 244  | 961459   | 68.485 | ng    | 0.00     |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

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

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