

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF020823\  
 Data File : BF132362.D  
 Acq On : 08 Feb 2023 17:59  
 Operator : CG\JU  
 Sample : 01418-04  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 JPP-16-020223

Quant Time: Feb 09 02:00:44 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF013123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 09 01:56:26 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.043  | 152  | 171707   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.331  | 136  | 651516   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 10.095 | 164  | 311627   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.589 | 188  | 503366   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.248 | 240  | 309579   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.824 | 264  | 224776   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.684  | 112  | 1280901  | 119.909 | ng    | 0.02     |
| 7) Phenol-d6                | 6.666  | 99   | 1395328  | 105.949 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.607  | 82   | 1023747  | 87.440  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.889 | 330  | 487989   | 152.275 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.413  | 172  | 1826797  | 90.526  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.189 | 244  | 1645567  | 102.823 | ng    | 0.00     |

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

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

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