

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF031422\  
 Data File : BF127317.D  
 Acq On : 14 Mar 2022 13:28  
 Operator : CG\JU  
 Sample : PB143209BL  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB143209BL

Quant Time: Mar 14 13:50:03 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF031022.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Mar 14 12:19:36 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.869  | 152  | 130313   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.151  | 136  | 484240   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.910  | 164  | 278234   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.398 | 188  | 505821   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 14.045 | 240  | 404563   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.527 | 264  | 316116   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.510  | 112  | 981795   | 121.116 | ng    | 0.02     |
| 7) Phenol-d6                | 6.504  | 99   | 1365476  | 122.449 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.434  | 82   | 966610   | 82.437  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.704 | 330  | 373752   | 124.785 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.228  | 172  | 1614827  | 82.054  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.992 | 244  | 1861236  | 79.220  | ng    | 0.00     |

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

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

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