

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF083023\  
 Data File : BF135028.D  
 Acq On : 31 Aug 2023 03:59  
 Operator : CG\JU  
 Sample : 04198-07  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OR-309-OK-370-COMP-25

Quant Time: Aug 31 06:01:46 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF083023.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Aug 31 03:48:08 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.769  | 152  | 122991   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.051  | 136  | 472820   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.810  | 164  | 210125   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.298 | 188  | 306222   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 13.957 | 240  | 236444   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.427 | 264  | 214435   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.410  | 112  | 880795   | 110.617 | ng    | 0.01     |
| 7) Phenol-d6                | 6.410  | 99   | 1072208  | 108.033 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.334  | 82   | 650868   | 77.336  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.598 | 330  | 208424   | 93.506  | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.133  | 172  | 1093474  | 84.725  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.898 | 244  | 864986   | 59.700  | ng    | 0.00     |

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

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

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