

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF070524\  
 Data File : BF138471.D  
 Acq On : 05 Jul 2024 16:33  
 Operator : CG\JU  
 Sample : P3135-01RE  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OK-628-COMP-01RE

Quant Time: Jul 05 17:03:52 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF062824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Jun 29 02:44:30 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.869  | 152  | 38144    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.151  | 136  | 144648   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.904  | 164  | 78441    | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.392 | 188  | 127191   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.027 | 240  | 85763    | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.498 | 264  | 62413    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.487  | 112  | 231089   | 100.444 | ng    | -0.01    |
| 7) Phenol-d6                | 6.498  | 99   | 300006   | 96.800  | ng    | -0.02    |
| 23) Nitrobenzene-d5         | 7.428  | 82   | 181794   | 65.288  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol    | 10.692 | 330  | 61611    | 96.136  | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.228  | 172  | 332410   | 67.441  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.974 | 244  | 297034   | 47.463  | ng    | 0.00     |

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

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

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