

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091924\  
 Data File : BF139483.D  
 Acq On : 20 Sep 2024 00:36  
 Operator : RC/JU  
 Sample : P4088-02  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 NB-614-COMP-07

Quant Time: Sep 20 01:42:08 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091324.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Sep 14 02:58:23 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.881  | 152  | 80791    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.157  | 136  | 297776   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.910  | 164  | 160031   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.398 | 188  | 255254   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.027 | 240  | 126749   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.498 | 264  | 112829   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.516  | 112  | 732200   | 147.847 | ng    | 0.02     |
| 7) Phenol-d6                | 6.510  | 99   | 919874   | 141.511 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.440  | 82   | 645392   | 101.891 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.704 | 330  | 228418   | 134.070 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.234  | 172  | 1149737  | 100.764 | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.980 | 244  | 816831   | 105.779 | ng    | 0.00     |

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

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

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