

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101924\  
Data File : BF139878.D  
Acq On : 19 Oct 2024 18:18  
Operator : RC/JU  
Sample : P4425-08  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-4

Quant Time: Oct 21 01:42:39 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101824.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Fri Oct 18 15:07:50 2024  
Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.892  | 152  | 205805   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.169  | 136  | 797916   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.927  | 164  | 434693   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.410 | 188  | 696781   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.045 | 240  | 383604   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.527 | 264  | 380773   | 20.000  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.516  | 112  | 1482632  | 112.779 | ng    | 0.02     |
| 7) Phenol-d6                | 6.510  | 99   | 1794080  | 105.369 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.451  | 82   | 1294650  | 89.927  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.716 | 330  | 544366   | 133.883 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.245  | 172  | 2213967  | 84.175  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.998 | 244  | 1964458  | 83.447  | ng    | 0.00     |

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

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

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