

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF103124\  
 Data File : BF140186.D  
 Acq On : 31 Oct 2024 13:55  
 Operator : RC/JU  
 Sample : P4611-06  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-2

Quant Time: Oct 31 14:29:53 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   | 6.886  | 152  | 93580    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.163  | 136  | 370111   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.922  | 164  | 216985   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.416 | 188  | 382650   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.092 | 240  | 213542   | 20.000  | ng    | 0.04     |
| 86) Perylene-d12            | 15.633 | 264  | 174560   | 20.000  | ng    | 0.11     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.510  | 112  | 799434   | 133.736 | ng    | 0.01     |
| 7) Phenol-d6                | 6.510  | 99   | 967049   | 124.908 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.445  | 82   | 762835   | 114.234 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.722 | 330  | 358347   | 176.560 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.239  | 172  | 1347457  | 102.631 | ng    | -0.01    |
| 79) Terphenyl-d14           | 13.015 | 244  | 1457276  | 111.201 | ng    | 0.02     |

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

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

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