

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF073124\  
 Data File : BF138710.D  
 Acq On : 31 Jul 2024 19:06  
 Operator : RC/JU  
 Sample : P3387-03  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 WC-9-Full Waste Class

Quant Time: Aug 01 04:33:36 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF073024.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jul 30 17:50:01 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.845  | 152  | 54464    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.128  | 136  | 192031   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.881  | 164  | 90712    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.363 | 188  | 166127   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.004 | 240  | 112809   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.468 | 264  | 86732    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.487  | 112  | 458527   | 129.959 | ng    | 0.02     |
| 7) Phenol-d6                | 6.487  | 99   | 563999   | 119.061 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.410  | 82   | 357102   | 90.919  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.669 | 330  | 111122   | 149.548 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.198  | 172  | 564164   | 93.445  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.951 | 244  | 713049   | 105.828 | ng    | 0.00     |

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

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

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