

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF042522\  
 Data File : BF127873.D  
 Acq On : 25 Apr 2022 11:44  
 Operator : CG\JU  
 Sample : N2487-03  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 2

Instrument :  
 BNA\_F  
 ClientSampleId :  
 UST-1-E

Quant Time: Apr 25 16:55:07 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF042222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sun Apr 24 00:43:03 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.940  | 152  | 278801   | 40.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.222  | 136  | 1082590  | 40.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.981  | 164  | 612972   | 40.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.475 | 188  | 1088280  | 40.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.121 | 240  | 708993   | 40.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.633 | 264  | 635519   | 40.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.569  | 112  | 1184467  | 135.337 | ng    | 0.00     |
| 7) Phenol-d6                | 6.569  | 99   | 1526523  | 134.421 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.504  | 82   | 1094134  | 95.051  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.775 | 330  | 398554   | 129.178 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.298  | 172  | 1666479  | 91.986  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.063 | 244  | 1655638  | 85.401  | ng    | 0.00     |

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

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

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