

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF122623\  
 Data File : BF136732.D  
 Acq On : 26 Dec 2023 22:05  
 Operator : CG\JU  
 Sample : 05933-03  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 S-2-NORTH-SIDE

Quant Time: Dec 27 03:59:56 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122023.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Dec 21 01:54:25 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.781  | 152  | 69087    | 20.000 | ng    | -0.01    |
| 21) Naphthalene-d8          | 8.057  | 136  | 261447   | 20.000 | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.816  | 164  | 123090   | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.304 | 188  | 193403   | 20.000 | ng    | -0.01    |
| 76) Chrysene-d12            | 13.962 | 240  | 133366   | 20.000 | ng    | # 0.00   |
| 86) Perylene-d12            | 15.439 | 264  | 124417   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.428  | 112  | 356712   | 80.561 | ng    | 0.00     |
| 7) Phenol-d6                | 6.428  | 99   | 451171   | 78.637 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.345  | 82   | 278099   | 54.430 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.610 | 330  | 81077    | 55.928 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.139  | 172  | 540963   | 59.110 | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.904 | 244  | 436744   | 45.764 | ng    | -0.01    |

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

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

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