

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF082421\  
 Data File : BF125192.D  
 Acq On : 24 Aug 2021 18:57  
 Operator : JU/CG  
 Sample : M3505-02  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 DRILL-CUTTING

Quant Time: Aug 25 04:36:33 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF081821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 24 12:39:21 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.916  | 152  | 148980   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.198  | 136  | 572793   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.951  | 164  | 312809   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.439 | 188  | 544001   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.080 | 240  | 375028   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.574 | 264  | 397140   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.539  | 112  | 1019425  | 103.571 | ng    | 0.01     |
| 7) Phenol-d6                | 6.539  | 99   | 1174727  | 92.281  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.481  | 82   | 925046   | 81.304  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.739 | 330  | 452363   | 144.871 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.274  | 172  | 1665939  | 74.226  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.027 | 244  | 1809184  | 76.853  | ng    | 0.00     |

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

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

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