

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061421\  
 Data File : BF124315.D  
 Acq On : 14 Jun 2021 20:00  
 Operator : JU/CG  
 Sample : M2625-34RE  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 18-B12(190-194)RE

Quant Time: Jun 15 02:27:30 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF060321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 14 12:10:02 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.828  | 152  | 32488    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.110  | 136  | 120146   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.869  | 164  | 28465    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.357 | 188  | 92026    | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 14.010 | 240  | 22343    | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.492 | 264  | 975      | 20.000  | ng    | # 0.00   |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.475  | 112  | 50766    | 23.523  | ng    | 0.02     |
| 7) Phenol-d6                | 6.487  | 99   | 27958    | 9.638   | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.392  | 82   | 111323   | 36.749  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.663 | 330  | 25321    | 62.788  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.186  | 172  | 210859   | 93.215  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.945 | 244  | 184270   | 113.232 | ng    | -0.01    |

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

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

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