

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110921\  
 Data File : BP007890.D  
 Acq On : 09 Nov 2021 14:16  
 Operator : CG/JU  
 Sample : M4576-03RE  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SB-2-(9-9.5)RE

Quant Time: Nov 09 14:44:13 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP110221.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 09 10:45:24 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.863  | 152  | 120531   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.663 | 136  | 407070   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.498 | 164  | 249093   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.251 | 188  | 526530   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.345 | 240  | 533319   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.757 | 264  | 595922   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.446  | 112  | 913498   | 128.534 | ng    | 0.00     |
| 7) Phenol-d6                | 7.040  | 99   | 1071401  | 104.154 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.016  | 82   | 629241   | 83.035  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 15.992 | 330  | 449613   | 126.476 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.122 | 172  | 1473779  | 86.162  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.816 | 244  | 2316118  | 88.433  | ng    | 0.00     |

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

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

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