

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP011022\  
 Data File : BP008745.D  
 Acq On : 10 Jan 2022 11:24  
 Operator : CG/JU  
 Sample : PB141870BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PB141870BL

Quant Time: Jan 10 16:33:22 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP010622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jan 07 07:45:52 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.869  | 152  | 349881   | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 10.675 | 136  | 1441209  | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.510 | 164  | 1010234  | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.263 | 188  | 2291479  | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.351 | 240  | 2451532  | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.774 | 264  | 2661769  | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.452  | 112  | 2851640  | 127.631 | ng    | 0.00     |
| 7) Phenol-d6                | 7.046  | 99   | 3571417  | 119.072 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.028  | 82   | 2193215  | 86.183  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.998 | 330  | 1951166  | 114.480 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 13.134 | 172  | 6170271  | 85.868  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.822 | 244  | 10862958 | 90.299  | ng    | 0.00     |

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

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

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