

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP021722\  
 Data File : BP009206.D  
 Acq On : 17 Feb 2022 21:14  
 Operator : CG/JU  
 Sample : N1561-02 5X  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TP-1-1-COMP

Quant Time: Feb 18 01:56:23 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP020722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Feb 16 14:30:24 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.775  | 152  | 275914   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.569 | 136  | 942110   | 20.000 | ng    | -0.01    |
| 39) Acenaphthene-d10        | 14.416 | 164  | 510659   | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10        | 17.181 | 188  | 986621   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 21.280 | 240  | 1166349  | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 23.668 | 264  | 1285673  | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.381  | 112  | 250720   | 16.209 | ng    | 0.00     |
| 7) Phenol-d6                | 6.981  | 99   | 273328   | 13.411 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.957  | 82   | 170052   | 10.179 | ng    | 0.01     |
| 42) 2,4,6-Tribromophenol    | 15.934 | 330  | 93942    | 13.778 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.045 | 172  | 422154   | 11.573 | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.739 | 244  | 595611   | 9.183  | ng    | -0.01    |

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

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

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