

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP071421\  
 Data File : BP006223.D  
 Acq On : 14 Jul 2021 17:52  
 Operator : CG/JU  
 Sample : M3026-06  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 ABN-JG-2

Quant Time: Jul 15 03:10:14 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP070721.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jul 12 14:59:51 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.852  | 152  | 283283   | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 10.652 | 136  | 1080830  | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 14.487 | 164  | 576040   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 17.240 | 188  | 1156046  | 20.000  | ng    | #-0.01   |
| 76) Chrysene-d12            | 21.316 | 240  | 1187360  | 20.000  | ng    | #-0.01   |
| 86) Perylene-d12            | 23.698 | 264  | 1342375  | 20.000  | ng    | -0.02    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.434  | 112  | 1910372  | 116.035 | ng    | 0.00     |
| 7) Phenol-d6                | 7.040  | 99   | 2076178  | 95.337  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.022  | 82   | 1374407  | 80.322  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 15.987 | 330  | 1111834  | 162.119 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 13.110 | 172  | 3164125  | 81.204  | ng    | -0.01    |
| 79) Terphenyl-d14           | 19.792 | 244  | 5539161  | 91.978  | ng    | -0.01    |

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

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

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