

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061821\  
 Data File : BP005972.D  
 Acq On : 18 Jun 2021 15:25  
 Operator : CG/JU  
 Sample : M2687-18  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DUP-01

Quant Time: Jun 18 16:15:51 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP060821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jun 16 11:35:30 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min)     |
|-----------------------------|--------|------|----------|---------|-------|--------------|
| -----                       |        |      |          |         |       |              |
| Internal Standards          |        |      |          |         |       |              |
| 1) 1,4-Dichlorobenzene-d4   | 7.928  | 152  | 66940    | 20.000  | ng    | 0.00         |
| 21) Naphthalene-d8          | 10.728 | 136  | 244858   | 20.000  | ng    | 0.00         |
| 39) Acenaphthene-d10        | 14.551 | 164  | 151554   | 20.000  | ng    | 0.00         |
| 64) Phenanthrene-d10        | 17.298 | 188  | 330392   | 20.000  | ng    | 0.00         |
| 76) Chrysene-d12            | 21.375 | 240  | 384000   | 20.000  | ng    | 0.00         |
| 86) Perylene-d12            | 23.780 | 264  | 441769   | 20.000  | ng    | -0.01        |
| System Monitoring Compounds |        |      |          |         |       |              |
| 5) 2-Fluorophenol           | 5.499  | 112  | 551050   | 132.093 | ng    | 0.00         |
| 7) Phenol-d6                | 7.105  | 99   | 571403   | 105.676 | ng    | 0.00         |
| 23) Nitrobenzene-d5         | 9.093  | 82   | 443724   | 88.845  | ng    | 0.00         |
| 42) 2,4,6-Tribromophenol    | 16.045 | 330  | 399383   | 153.563 | ng    | 0.00         |
| 45) 2-Fluorobiphenyl        | 13.181 | 172  | 994931   | 86.205  | ng    | 0.00         |
| 79) Terphenyl-d14           | 19.851 | 244  | 1929726  | 94.099  | ng    | 0.00         |
| Target Compounds            |        |      |          |         |       |              |
| 32) Benzoic acid            | 10.016 | 122  | 2342     | 7.056   | ng    | Qvalue<br>88 |
| -----                       |        |      |          |         |       |              |

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

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