

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_E\DATA\BE101618\  
 Data File : BE097941.D  
 Acq On : 17 Oct 2018 10:53  
 Operator : SJ/JU  
 Sample : J5393-10DL 2X  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 RE137-700FT-20181009DL

Quant Time: Oct 17 17:09:38 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_E\METHODS\8270-SIM-BE101618.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Oct 16 15:36:44 2018  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units | Dev(Min)  |
|-----------------------------|-------|------|----------|-------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.88  | 152  | 1922     | 0.40  | ng    | 0.00      |
| 7) Naphthalene-d8           | 10.68 | 136  | 7943     | 0.40  | ng    | -0.01     |
| 13) Acenaphthene-d10        | 14.54 | 164  | 5051     | 0.40  | ng    | -0.02     |
| 19) Phenanthrene-d10        | 17.29 | 188  | 13965    | 0.40  | ng    | -0.01     |
| 27) Chrysene-d12            | 21.47 | 240  | 17463    | 0.40  | ng    | -0.01     |
| 34) Perylene-d12            | 24.03 | 264  | 18171    | 0.40  | ng    | -0.01     |
| System Monitoring Compounds |       |      |          |       |       |           |
| 4) 2-Fluorophenol           | 5.46  | 112  | 547      | 0.09  | ng    | 0.00      |
| 5) Phenol-d6                | 7.05  | 99   | 451      | 0.05  | ng    | 0.00      |
| 8) Nitrobenzene-d5          | 9.04  | 82   | 1284     | 0.17  | ng    | 0.00      |
| 11) 2-Methylnaphthalene-d10 | 12.28 | 152  | 2278     | 0.17  | ng    | -0.02     |
| 14) 2,4,6-Tribromophenol    | 16.04 | 330  | 417      | 0.15  | ng    | 0.00      |
| 15) 2-Fluorobiphenyl        | 13.17 | 172  | 4064     | 0.20  | ng    | 0.00      |
| 25) Fluoranthene-d10        | 19.33 | 212  | 33835    | 0.18  | ng    | 0.00      |
| 29) Terphenyl-d14           | 19.92 | 244  | 11061    | 0.28  | ng    | -0.01     |
| Target Compounds            |       |      |          |       |       |           |
| 2) 1,4-Dioxane              | 3.37  | 88   | 9266     | 2.765 | ng    | Qvalue 94 |
| 6) bis(2-Chloroethyl)ether  | 7.30  | 93   | 212      | 0.028 | ng    | # 83      |

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

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