

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121318\  
 Data File : BE098402.D  
 Acq On : 13 Dec 2018 18:22  
 Operator : SJ/JU  
 Sample : J6323-02  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 RE-104D2-20181206

Quant Time: Dec 13 21:09:46 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Dec 10 14:59:55 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.86  | 152  | 888      | 0.40 | ng    | 0.00     |
| 7) Naphthalene-d8         | 10.67 | 136  | 3976     | 0.40 | ng    | 0.00     |
| 13) Acenaphthene-d10      | 14.53 | 164  | 2666     | 0.40 | ng    | 0.00     |
| 19) Phenanthrene-d10      | 17.27 | 188  | 7286     | 0.40 | ng    | -0.01    |
| 27) Chrysene-d12          | 21.46 | 240  | 8679     | 0.40 | ng    | 0.00     |
| 34) Perylene-d12          | 24.00 | 264  | 8281     | 0.40 | ng    | -0.01    |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|-------|------|----------|------|-------|----------|
| 4) 2-Fluorophenol           | 5.46  | 112  | 453      | 0.22 | ng    | 0.01     |
| 5) Phenol-d6                | 7.03  | 99   | 388      | 0.13 | ng    | -0.01    |
| 8) Nitrobenzene-d5          | 9.03  | 82   | 1350     | 0.37 | ng    | 0.01     |
| 11) 2-Methylnaphthalene-d10 | 12.27 | 152  | 2591     | 0.40 | ng    | 0.00     |
| 14) 2,4,6-Tribromophenol    | 16.02 | 330  | 589      | 0.60 | ng    | 0.00     |
| 15) 2-Fluorobiphenyl        | 13.14 | 172  | 4767     | 0.42 | ng    | 0.00     |
| 25) Fluoranthene-d10        | 19.30 | 212  | 41206    | 0.44 | ng    | 0.00     |
| 29) Terphenyl-d14           | 19.90 | 244  | 8848     | 0.42 | ng    | 0.00     |

| Target Compounds               | R.T.  | QIon | Response | Conc  | Units | Qvalue |
|--------------------------------|-------|------|----------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.37  | 88   | 654      | 0.426 | ng    | # 50   |
| 6) bis(2-Chloroethyl)ether     | 7.29  | 93   | 328      | 0.117 | ng    | # 71   |
| 23) Phenanthrene               | 17.32 | 178  | 479      | 0.027 | ng    | # 85   |
| 26) Fluoranthene               | 19.34 | 202  | 728      | 0.031 | ng    | # 96   |
| 28) Pyrene                     | 19.70 | 202  | 671      | 0.025 | ng    | # 99   |
| 30) Benzo(a)anthracene         | 21.44 | 228  | 646      | 0.026 | ng    | # 80   |
| 31) Chrysene                   | 21.50 | 228  | 645      | 0.025 | ng    | # 74   |
| 32) Bis(2-ethylhexyl)phthalate | 21.35 | 149  | 4224     | 0.084 | ng    | # 100  |
| 33) Indeno(1,2,3-cd)pyrene     | 26.71 | 276  | 524      | 0.020 | ng    | # 76   |

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

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