

Data Path : Z:\HPCHEM1\BNA E\DATA\BE043018\  
 Data File : BE096224.D  
 Acq On : 30 Apr 2018 16:33  
 Operator : SJ/JU  
 Sample : J2463-12MSD  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 LF001MW001-180418MSD

Quant Time: May 01 02:28:29 2018  
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE041118.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Apr 30 14:27:29 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.36  | 152  | 807      | 0.40 | ng    | 0.00     |
| 7) Naphthalene-d8         | 11.20 | 136  | 2722     | 0.40 | ng    | 0.01     |
| 13) Acenaphthene-d10      | 14.93 | 164  | 8712     | 0.40 | ng    | -0.01    |
| 19) Phenanthrene-d10      | 17.65 | 188  | 3578     | 0.40 | ng    | 0.00     |
| 27) Chrysene-d12          | 21.75 | 240  | 3700     | 0.40 | ng    | 0.01     |
| 34) Perylene-d12          | 24.35 | 264  | 3596     | 0.40 | ng    | 0.00     |

|                             |       |     |       |      |    |      |
|-----------------------------|-------|-----|-------|------|----|------|
| System Monitoring Compounds |       |     |       |      |    |      |
| 4) 2-Fluorophenol           | 5.88  | 112 | 618   | 0.25 | ng | 0.01 |
| 5) Phenol-d6                | 7.53  | 99  | 818   | 0.24 | ng | 0.02 |
| 8) Nitrobenzene-d5          | 9.54  | 82  | 2525  | 0.79 | ng | 0.00 |
| 11) 2-Methylnaphthalene-d10 | 12.74 | 152 | 1653  | 0.38 | ng | 0.00 |
| 14) 2,4,6-Tribromophenol    | 16.41 | 330 | 320   | 0.08 | ng | 0.00 |
| 15) 2-Fluorobiphenyl        | 13.59 | 172 | 2395  | 0.06 | ng | 0.01 |
| 25) Fluoranthene-d10        | 19.64 | 212 | 16626 | 0.34 | ng | 0.00 |
| 29) Terphenyl-d14           | 20.19 | 244 | 2563  | 0.31 | ng | 0.00 |

|                                |       |     |         |         |    |      |
|--------------------------------|-------|-----|---------|---------|----|------|
| Target Compounds               |       |     |         | Qvalue  |    |      |
| 2) 1,4-Dioxane                 | 3.61  | 88  | 437     | 0.351   | ng | # 1  |
| 3) n-Nitrosodimethylamine      | 3.99  | 42  | 642370  | 420.605 | ng | # 1  |
| 6) bis(2-Chloroethyl)ether     | 7.76  | 93  | 1103    | 0.364   | ng | # 1  |
| 9) Naphthalene                 | 11.26 | 128 | 3253495 | 113.378 | ng | 99   |
| 10) Hexachlorobutadiene        | 11.50 | 225 | 176     | 0.138   | ng | 90   |
| 12) 2-Methylnaphthalene        | 12.81 | 142 | 81413   | 17.140  | ng | 94   |
| 16) Acenaphthylene             | 14.67 | 152 | 3746    | 0.081   | ng | # 91 |
| 17) Acenaphthene               | 15.00 | 154 | 2159    | 0.077   | ng | # 92 |
| 18) Fluorene                   | 15.97 | 166 | 4034    | 0.117   | ng | # 81 |
| 20) 4-Bromophenyl-phenylether  | 16.85 | 248 | 755     | 0.355   | ng | # 38 |
| 21) Hexachlorobenzene          | 16.97 | 284 | 684     | 0.323   | ng | # 77 |
| 22) Pentachlorophenol          | 17.31 | 266 | 920     | 1.077   | ng | # 75 |
| 23) Phenanthrene               | 17.70 | 178 | 4515    | 0.449   | ng | 97   |
| 24) Anthracene                 | 17.78 | 178 | 3858    | 0.402   | ng | # 92 |
| 26) Fluoranthene               | 19.66 | 202 | 5249    | 0.397   | ng | # 25 |
| 28) Pyrene                     | 19.97 | 202 | 541     | 0.074   | ng | # 78 |
| 30) Benzo(a)anthracene         | 21.73 | 228 | 4399    | 0.371   | ng | # 92 |
| 31) Chrysene                   | 21.79 | 228 | 3928    | 0.343   | ng | # 87 |
| 32) Bis(2-ethylhexyl)phthalate | 21.60 | 149 | 12003   | 0.392   | ng | 99   |
| 33) Indeno(1,2,3-cd)pyrene     | 27.13 | 276 | 2708    | 0.244   | ng | 95   |
| 35) Benzo(b)fluoranthene       | 23.55 | 252 | 3302    | 0.306   | ng | # 87 |
| 36) Benzo(k)fluoranthene       | 23.60 | 252 | 3112    | 0.280   | ng | # 67 |
| 37) Benzo(a)pyrene             | 24.24 | 252 | 2980    | 0.287   | ng | # 71 |
| 38) Dibenzo(a,h)anthracene     | 27.15 | 278 | 2203    | 0.241   | ng | # 76 |
| 39) Benzo(g,h,i)perylene       | 28.00 | 276 | 2294    | 0.237   | ng | # 83 |

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

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