

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE030618\  
 Data File : BE095704.D  
 Acq On : 6 Mar 2018 19:11  
 Operator : SJ/JU  
 Sample : J1811-01MS 20X  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 T-1\_030518MS

Manual Integrations  
 APPROVED

Sohil  
 3/7/2018 12:20:10 PM

Quant Time: Mar 07 01:46:36 2018  
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE022318.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Feb 23 12:12:42 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.43  | 152  | 1378     | 0.40 | ng    | 0.00     |
| 7) Naphthalene-d8         | 11.27 | 136  | 5570     | 0.40 | ng    | 0.00     |
| 13) Acenaphthene-d10      | 15.00 | 164  | 3232     | 0.40 | ng    | 0.00     |
| 19) Phenanthrene-d10      | 17.71 | 188  | 7380     | 0.40 | ng    | 0.00     |
| 27) Chrysene-d12          | 21.80 | 240  | 7591     | 0.40 | ng    | 0.00     |
| 34) Perylene-d12          | 24.43 | 264  | 7118     | 0.40 | ng    | 0.00     |

|                             |       |     |     |      |    |      |
|-----------------------------|-------|-----|-----|------|----|------|
| System Monitoring Compounds |       |     |     |      |    |      |
| 4) 2-Fluorophenol           | 5.92  | 112 | 43  | 0.01 | ng | 0.00 |
| 5) Phenol-d6                | 7.57  | 99  | 57  | 0.01 | ng | 0.01 |
| 8) Nitrobenzene-d5          | 9.61  | 82  | 79  | 0.01 | ng | 0.00 |
| 11) 2-Methylnaphthalene-d10 | 12.81 | 152 | 123 | 0.01 | ng | 0.00 |
| 14) 2,4,6-Tribromophenol    | 16.47 | 330 | 17  | 0.13 | ng | 0.00 |
| 15) 2-Fluorobiphenyl        | 13.64 | 172 | 192 | 0.01 | ng | 0.00 |
| 25) Fluoranthene-d10        | 19.69 | 212 | 993 | 0.01 | ng | 0.00 |
| 29) Terphenyl-d14           | 20.25 | 244 | 191 | 0.01 | ng | 0.00 |

|                                |       |     |       |        |    |      |
|--------------------------------|-------|-----|-------|--------|----|------|
| Target Compounds               |       |     |       | Qvalue |    |      |
| 2) 1,4-Dioxane                 | 3.64  | 88  | 38    | 0.018  | ng | # 82 |
| 3) n-Nitrosodimethylamine      | 4.01  | 42  | 44    | 0.016  | ng | # 3  |
| 6) bis(2-Chloroethyl)ether     | 7.83  | 93  | 64    | 0.012  | ng | # 96 |
| 9) Naphthalene                 | 11.31 | 128 | 1419  | 0.022  | ng | # 86 |
| 10) Hexachlorobutadiene        | 11.58 | 225 | 52    | 0.014  | ng | # 94 |
| 12) 2-Methylnaphthalene        | 12.88 | 142 | 571   | 0.052  | ng | # 94 |
| 16) Acenaphthylene             | 14.74 | 152 | 1007  | 0.055  | ng | # 92 |
| 17) Acenaphthene               | 15.07 | 154 | 187   | 0.016  | ng | # 75 |
| 18) Fluorene                   | 16.03 | 166 | 585   | 0.041  | ng | # 74 |
| 20) 4-Bromophenyl-phenylether  | 16.90 | 248 | 66    | 0.014  | ng | # 60 |
| 21) Hexachlorobenzene          | 17.02 | 284 | 55    | 0.011  | ng | # 65 |
| 22) Pentachlorophenol          | 17.36 | 266 | 24    | 0.192  | ng | # 97 |
| 23) Phenanthrene               | 17.75 | 178 | 13573 | 0.588  | ng | # 99 |
| 24) Anthracene                 | 17.84 | 178 | 2424  | 0.114  | ng | # 93 |
| 26) Fluoranthene               | 19.72 | 202 | 23100 | 0.814  | ng | # 97 |
| 28) Pyrene                     | 20.07 | 202 | 26941 | 0.868  | ng | # 97 |
| 30) Benzo(a)anthracene         | 21.78 | 228 | 9729  | 0.393  | ng | # 97 |
| 31) Chrysene                   | 21.84 | 228 | 9455  | 0.384  | ng | # 95 |
| 32) Bis(2-ethylhexyl)phthalate | 21.66 | 149 | 1357  | 0.031  | ng | # 96 |
| 33) Indeno(1,2,3-cd)pyrene     | 27.23 | 276 | 4246  | 0.178  | ng | # 91 |
| 35) Benzo(b)fluoranthene       | 23.61 | 252 | 9271  | 0.366  | ng | # 92 |
| 36) Benzo(k)fluoranthene       | 23.66 | 252 | 3799m | 0.151  | ng | # 92 |
| 37) Benzo(a)pyrene             | 24.30 | 252 | 6426  | 0.281  | ng | # 92 |
| 38) Dibenzo(a,h)anthracene     | 27.24 | 278 | 1274  | 0.063  | ng | # 75 |
| 39) Benzo(g,h,i)perylene       | 28.10 | 276 | 5048  | 0.232  | ng | # 89 |

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

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