

Data Path : Z:\HPCHEM1\BNA E\DATA\BE110816\  
 Data File : BE092512.D  
 Acq On : 8 Nov 2016 17:07  
 Operator : UM/SJ  
 Sample : PB94535BSD  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 PB94535BSD

Manual Integrations  
 APPROVED

umangi  
 11/9/2016 10:58:55 AM

Quant Time: Nov 09 00:52:30 2016  
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE110816.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 08 16:27:22 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.14  | 152  | 9032     | 5.00 | ng    | -0.02    |
| 7) Naphthalene-d8         | 10.93 | 136  | 40708    | 5.00 | ng    | 0.00     |
| 12) Acenaphthene-d10      | 14.80 | 164  | 26544    | 5.00 | ng    | 0.00     |
| 18) Phenanthrene-d10      | 17.56 | 188  | 50768    | 5.00 | ng    | 0.00     |
| 24) Chrysene-d12          | 21.72 | 240  | 64045    | 5.00 | ng    | 0.00     |
| 31) Perylene-d12          | 24.09 | 264  | 56694    | 5.00 | ng    | 0.00     |

|                             |       |     |       |       |    |       |
|-----------------------------|-------|-----|-------|-------|----|-------|
| System Monitoring Compounds |       |     |       |       |    |       |
| 4) 2-Fluorophenol           | 5.65  | 112 | 20507 | 10.30 | ng | -0.01 |
| 5) Phenol-d6                | 7.31  | 99  | 28653 | 10.79 | ng | -0.02 |
| 8) Nitrobenzene-d5          | 9.32  | 82  | 13423 | 5.03  | ng | -0.02 |
| 13) 2,4,6-Tribromophenol    | 16.30 | 330 | 7176  | 7.58  | ng | -0.01 |
| 14) 2-Fluorobiphenyl        | 13.42 | 172 | 30043 | 4.58  | ng | -0.01 |
| 26) Terphenyl-d14           | 20.16 | 244 | 38766 | 4.98  | ng | -0.02 |

|                                |       |     |         |       |    |        |
|--------------------------------|-------|-----|---------|-------|----|--------|
| Target Compounds               |       |     |         |       |    | Qvalue |
| 2) 1,4-Dioxane                 | 3.39  | 88  | 4265    | 3.98  | ng | # 55   |
| 3) n-Nitrosodimethylamine      | 3.73  | 42  | 6783    | 4.24  | ng | # 97   |
| 6) bis(2-Chloroethyl)ether     | 7.57  | 93  | 11614   | 3.99  | ng | 90     |
| 9) Naphthalene                 | 10.97 | 128 | 36419   | 4.14  | ng | # 95   |
| 10) Hexachlorobutadiene        | 11.26 | 225 | 5554    | 4.16  | ng | 100    |
| 11) 2-Methylnaphthalene        | 12.60 | 142 | 24626   | 4.30  | ng | 97     |
| 15) Acenaphthylene             | 14.51 | 152 | 162688  | 4.27  | ng | 100    |
| 16) Acenaphthene               | 14.87 | 154 | 24148   | 3.92  | ng | 91     |
| 17) Fluorene                   | 15.84 | 166 | 32293   | 4.18  | ng | 99     |
| 19) Hexachlorobenzene          | 16.87 | 284 | 9597    | 5.01  | ng | # 67   |
| 20) Pentachlorophenol          | 17.21 | 266 | 7872    | 14.21 | ng | # 85   |
| 21) Phenanthrene               | 17.58 | 178 | 48575   | 4.43  | ng | # 94   |
| 22) Anthracene                 | 17.67 | 178 | 51631   | 4.85  | ng | 99     |
| 23) Fluoranthene               | 19.59 | 202 | 234067  | 4.25  | ng | 98     |
| 25) Pyrene                     | 19.95 | 202 | 245670  | 4.68  | ng | 99     |
| 27) Benzo(a)anthracene         | 21.70 | 228 | 280258m | 4.32  | ng |        |
| 28) Chrysene                   | 21.75 | 228 | 241143m | 4.66  | ng |        |
| 29) Bis(2-ethylhexyl)phthalate | 21.61 | 149 | 22896   | 4.03  | ng | # 56   |
| 30) Indeno(1,2,3-cd)pyrene     | 26.55 | 276 | 312143  | 4.41  | ng | 98     |
| 32) Benzo(b)fluoranthene       | 23.35 | 252 | 65347   | 4.49  | ng | 95     |
| 33) Benzo(k)fluoranthene       | 23.39 | 252 | 70452   | 4.59  | ng | 93     |
| 34) Benzo(a)pyrene             | 23.97 | 252 | 62510   | 4.27  | ng | # 94   |
| 35) Dibenzo(a,h)anthracene     | 26.58 | 278 | 64978   | 4.71  | ng | 94     |
| 36) Benzo(g,h,i)perylene       | 27.31 | 276 | 270561  | 4.76  | ng | 98     |

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

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