

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN032119\  
 Data File : BN004853.D  
 Acq On : 21 Mar 2019 17:01  
 Operator : JU/SJ  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTDCCC0.4EC

Manual Integrations  
 APPROVED

Sohil  
 3/22/2019 7:16:20 PM

Quant Time: Mar 22 02:30:32 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN032019.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Mar 20 16:06:20 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.70  | 152  | 1899     | 0.40 | ng    | 0.00     |
| 7) Naphthalene-d8         | 10.48 | 136  | 8640     | 0.40 | ng    | -0.01    |
| 13) Acenaphthene-d10      | 14.33 | 164  | 5560     | 0.40 | ng    | -0.01    |
| 19) Phenanthrene-d10      | 17.09 | 188  | 14093    | 0.40 | ng    | 0.00     |
| 27) Chrysene-d12          | 21.29 | 240  | 18964    | 0.40 | ng    | 0.00     |
| 34) Perylene-d12          | 23.54 | 264  | 20894    | 0.40 | ng    | 0.00     |

|                             |       |     |       |      |    |      |
|-----------------------------|-------|-----|-------|------|----|------|
| System Monitoring Compounds |       |     |       |      |    |      |
| 4) 2-Fluorophenol           | 5.29  | 112 | 1985  | 0.38 | ng | 0.00 |
| 5) Phenol-d6                | 6.88  | 99  | 2541  | 0.40 | ng | 0.00 |
| 8) Nitrobenzene-d5          | 8.87  | 82  | 2141  | 0.37 | ng | 0.00 |
| 11) 2-Methylnaphthalene-d10 | 12.08 | 152 | 6018  | 0.40 | ng | 0.00 |
| 14) 2,4,6-Tribromophenol    | 15.85 | 330 | 1325  | 0.35 | ng | 0.00 |
| 15) 2-Fluorobiphenyl        | 12.96 | 172 | 9363  | 0.41 | ng | 0.00 |
| 25) Fluoranthene-d10        | 19.13 | 212 | 18091 | 0.38 | ng | 0.00 |
| 29) Terphenyl-d14           | 19.73 | 244 | 15309 | 0.39 | ng | 0.00 |

|                                |       |     |       |       |        |       |
|--------------------------------|-------|-----|-------|-------|--------|-------|
| Target Compounds               |       |     |       |       | Qvalue |       |
| 2) 1,4-Dioxane                 | 3.24  | 88  | 836   | 0.387 | ng     | 99    |
| 3) n-Nitrosodimethylamine      | 3.59  | 42  | 446m  | 0.350 | ng     |       |
| 6) bis(2-Chloroethyl)ether     | 7.14  | 93  | 2294  | 0.396 | ng     | 98    |
| 9) Naphthalene                 | 10.53 | 128 | 9736  | 0.402 | ng     | 99    |
| 10) Hexachlorobutadiene        | 10.80 | 225 | 1833  | 0.397 | ng     | # 100 |
| 12) 2-Methylnaphthalene        | 12.15 | 142 | 7244  | 0.402 | ng     | 99    |
| 16) Acenaphthylene             | 14.05 | 152 | 11317 | 0.395 | ng     | 100   |
| 17) Acenaphthene               | 14.40 | 154 | 7638  | 0.407 | ng     | 99    |
| 18) Fluorene                   | 15.39 | 166 | 10474 | 0.400 | ng     | 100   |
| 20) 4-Bromophenyl-phenylether  | 16.28 | 248 | 3606  | 0.392 | ng     | 95    |
| 21) Hexachlorobenzene          | 16.39 | 284 | 4264  | 0.405 | ng     | 99    |
| 22) Pentachlorophenol          | 16.77 | 266 | 421m  | 0.204 | ng     |       |
| 23) Phenanthrene               | 17.13 | 178 | 18254 | 0.401 | ng     | 99    |
| 24) Anthracene                 | 17.23 | 178 | 15813 | 0.389 | ng     | 100   |
| 26) Fluoranthene               | 19.16 | 202 | 22876 | 0.391 | ng     | 100   |
| 28) Pyrene                     | 19.52 | 202 | 23372 | 0.404 | ng     | 100   |
| 30) Benzo(a)anthracene         | 21.28 | 228 | 24226 | 0.390 | ng     | 100   |
| 31) Chrysene                   | 21.33 | 228 | 26717 | 0.408 | ng     | 99    |
| 32) Bis(2-ethylhexyl)phthalate | 21.18 | 149 | 9438  | 0.380 | ng     | 100   |
| 33) Indeno(1,2,3-cd)pyrene     | 25.81 | 276 | 36909 | 0.389 | ng     | 100   |
| 35) Benzo(b)fluoranthene       | 22.86 | 252 | 30544 | 0.404 | ng     | 99    |
| 36) Benzo(k)fluoranthene       | 22.91 | 252 | 29924 | 0.402 | ng     | 99    |
| 37) Benzo(a)pyrene             | 23.45 | 252 | 27712 | 0.400 | ng     | 99    |
| 38) Dibenzo(a,h)anthracene     | 25.82 | 278 | 30316 | 0.377 | ng     | 99    |
| 39) Benzo(g,h,i)perylene       | 26.52 | 276 | 30197 | 0.369 | ng     | 99    |

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN032119\  
Data File : BN004853.D  
Acq On : 21 Mar 2019 17:01  
Operator : JU/SJ  
Sample : SSTDCCC0.4  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTDCCC0.4EC

Manual Integrations  
APPROVED

Sohil  
3/22/2019 7:16:20 PM

Quant Time: Mar 22 02:30:32 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN032019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Wed Mar 20 16:06:20 2019  
Response via : Initial Calibration

