

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN020421\  
 Data File : BN013579.D  
 Acq On : 04 Feb 2021 19:42  
 Operator : CG/JU  
 Sample : SSTDC00.4EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD0.411

Quant Time: Feb 04 23:22:56 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN020421.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 04 14:51:05 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.56  | 152  | 4746     | 0.40 | ng/ul | 0.00     |
| 2) Naphthalene-d8         | 10.31 | 136  | 17111    | 0.40 | ng/ul | 0.00     |
| 6) Acenaphthene-d10       | 14.19 | 164  | 9107     | 0.40 | ng/ul | 0.00     |
| 10) Phenanthrene-d10      | 16.94 | 188  | 18945    | 0.40 | ng/ul | 0.00     |
| 16) Chrysene-d12          | 21.14 | 240  | 17459    | 0.40 | ng/ul | 0.00     |
| 20) Perylene-d12          | 23.31 | 264  | 14656    | 0.40 | ng/ul | 0.00     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|-------|------|----------|------|-------|----------|
| 4) 2-Methylnaphthalene-d10  | 11.91 | 152  | 10196    | 0.42 | ng/ul | 0.00     |
| 14) Fluoranthene-d10        | 18.98 | 212  | 19324    | 0.42 | ng/ul | 0.00     |

| Target Compounds           | R.T.  | QIon | Response | Conc  | Units  | Ovalue |
|----------------------------|-------|------|----------|-------|--------|--------|
| 3) Naphthalene             | 10.36 | 128  | 18759    | 0.421 | ng/ul  | 99     |
| 5) 2-Methylnaphthalene     | 11.98 | 142  | 12426    | 0.416 | ng/ul  | 99     |
| 7) Acenaphthylene          | 13.90 | 152  | 14117    | 0.416 | ng/ul  | 99     |
| 8) Acenaphthene            | 14.25 | 153  | 12460    | 0.413 | ng/ul  | 99     |
| 9) Fluorene                | 15.24 | 166  | 13945    | 0.408 | ng/ul  | 99     |
| 11) Pentachlorophenol      | 16.59 | 266  | 1032     | 0.447 | ng/ul  | 97     |
| 12) Phenanthrene           | 16.98 | 178  | 22956    | 0.407 | ng/ul  | 100    |
| 13) Anthracene             | 17.07 | 178  | 19190    | 0.410 | ng/ul  | 99     |
| 15) Fluoranthene           | 19.00 | 202  | 24385    | 0.401 | ng/ul  | 98     |
| 17) Pyrene                 | 19.37 | 202  | 24768    | 0.395 | ng/ul  | 100    |
| 18) Benzo(a)anthracene     | 21.13 | 228  | 22952    | 0.405 | ng/ul  | 99     |
| 19) Chrysene               | 21.18 | 228  | 25553    | 0.414 | ng/ul  | 100    |
| 21) Benzo(b)fluoranthene   | 22.66 | 252  | 23117    | 0.386 | ng/ul  | 99     |
| 22) Benzo(k)fluoranthene   | 22.70 | 252  | 25447    | 0.410 | ng/ul  | 99     |
| 23) Benzo(a)pyrene         | 23.21 | 252  | 19758    | 0.410 | ng/ul  | 98     |
| 24) Indeno(1,2,3-cd)pyrene | 25.46 | 276  | 26279    | 0.410 | ng/ul# | 100    |
| 25) Dibenzo(a,h)anthracene | 25.47 | 278  | 21266    | 0.404 | ng/ul  | 99     |
| 26) Benzo(g,h,i)perylene   | 26.11 | 276  | 22381    | 0.398 | ng/ul  | 99     |

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

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