

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN121718\  
 Data File : BN004066.D  
 Acq On : 18 Dec 2018 09:04  
 Operator : JU/SJ  
 Sample : J6270-14  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 F9E13

Manual Integrations  
 APPROVED

Sohil  
 12/18/2018 5:57:23 PM

Quant Time: Dec 18 11:33:56 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-SIM-BN111918.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Dec 18 06:35:27 2018  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.84  | 152  | 1614     | 0.40   | ng/ul  | 0.00     |
| 2) Naphthalene-d8           | 10.63 | 136  | 8250     | 0.40   | ng/ul  | 0.00     |
| 6) Acenaphthene-d10         | 14.47 | 164  | 5257     | 0.40   | ng/ul  | 0.00     |
| 10) Phenanthrene-d10        | 17.22 | 188  | 14054m   | 0.40   | ng/ul  | 0.00     |
| 16) Chrysene-d12            | 21.40 | 240  | 16033m   | 0.40   | ng/ul  | 0.00     |
| 20) Perylene-d12            | 23.72 | 264  | 13788m   | 0.40   | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |        |        |          |
| 4) 2-Methylnaphthalene-d10  | 12.22 | 152  | 2632     | 0.21   | ng/ul  | 0.00     |
| 14) Fluoranthene-d10        | 19.24 | 212  | 9624m    | 0.23   | ng/ul  | 0.00     |
| Target Compounds            |       |      |          | Qvalue |        |          |
| 7) Acenaphthylene           | 14.19 | 152  | 3427     | 0.145  | ng/ul  | 96       |
| 12) Phenanthrene            | 17.26 | 178  | 2161m    | 0.047  | ng/ul  |          |
| 13) Anthracene              | 17.35 | 178  | 3482m    | 0.092  | ng/ul  |          |
| 15) Fluoranthene            | 19.28 | 202  | 19551m   | 0.359  | ng/ul  |          |
| 17) Pyrene                  | 19.64 | 202  | 20683    | 0.360  | ng/ul  | 98       |
| 18) Benzo(a)anthracene      | 21.39 | 228  | 12789    | 0.253  | ng/ul  | 93       |
| 19) Chrysene                | 21.44 | 228  | 14294m   | 0.247  | ng/ul  |          |
| 21) Benzo(b)fluoranthene    | 23.02 | 252  | 22539m   | 0.395  | ng/ul  |          |
| 22) Benzo(k)fluoranthene    | 23.06 | 252  | 7635m    | 0.137  | ng/ul  |          |
| 23) Benzo(a)pyrene          | 23.62 | 252  | 9941     | 0.201  | ng/ul# | 82       |
| 24) Indeno(1,2,3-cd)pyrene  | 26.09 | 276  | 9376     | 0.166  | ng/ul# | 96       |
| 25) Dibenzo(a,h)anthracene  | 26.09 | 278  | 2387     | 0.052  | ng/ul# | 28       |
| 26) Benzo(g,h,i)perylene    | 26.82 | 276  | 7364     | 0.150  | ng/ul# | 81       |

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

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-SIM-BN11918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Tue Dec 18 06:35:27 2018  
Response via : Initial Calibration

