

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\  
 Data File : BN003973.D  
 Acq On : 15 Dec 2018 23:57  
 Operator : JU/SJ  
 Sample : J6171-07DL 5X  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 F9Y11DL

Manual Integrations  
 APPROVED

Sohil  
 12/17/2018 4:24:02 PM

Quant Time: Dec 16 02:10:24 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 : Sat Dec 15 23:04:58 2018  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.84  | 152  | 2000     | 0.40   | ng/ul  | 0.00     |
| 2) Naphthalene-d8           | 10.63 | 136  | 10145    | 0.40   | ng/ul  | 0.00     |
| 6) Acenaphthene-d10         | 14.47 | 164  | 6037     | 0.40   | ng/ul  | 0.00     |
| 10) Phenanthrene-d10        | 17.22 | 188  | 15070    | 0.40   | ng/ul  | 0.00     |
| 16) Chrysene-d12            | 21.40 | 240  | 19301    | 0.40   | ng/ul  | 0.00     |
| 20) Perylene-d12            | 23.71 | 264  | 17416    | 0.40   | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |        |        |          |
| 4) 2-Methylnaphthalene-d10  | 12.22 | 152  | 622      | 0.04   | ng/ul  | 0.00     |
| 14) Fluoranthene-d10        | 19.24 | 212  | 1810     | 0.04   | ng/ul  | 0.00     |
| Target Compounds            |       |      |          | Ovalue |        |          |
| 5) 2-Methylnaphthalene      | 12.29 | 142  | 472      | 0.024  | ng/ul  | 99       |
| 7) Acenaphthylene           | 14.19 | 152  | 619      | 0.023  | ng/ul# | 69       |
| 12) Phenanthrene            | 17.26 | 178  | 3058     | 0.062  | ng/ul# | 93       |
| 13) Anthracene              | 17.35 | 178  | 1117     | 0.027  | ng/ul# | 79       |
| 15) Fluoranthene            | 19.27 | 202  | 4322     | 0.074  | ng/ul  | 91       |
| 17) Pyrene                  | 19.63 | 202  | 5874     | 0.085  | ng/ul# | 93       |
| 18) Benzo(a)anthracene      | 21.38 | 228  | 3230     | 0.053  | ng/ul# | 73       |
| 19) Chrysene                | 21.43 | 228  | 5292     | 0.076  | ng/ul# | 79       |
| 21) Benzo(b)fluoranthene    | 23.01 | 252  | 6896m    | 0.096  | ng/ul  |          |
| 22) Benzo(k)fluoranthene    | 23.06 | 252  | 2365m    | 0.034  | ng/ul  |          |
| 23) Benzo(a)pyrene          | 23.61 | 252  | 3006     | 0.048  | ng/ul# | 10       |
| 24) Indeno(1,2,3-cd)pyrene  | 26.07 | 276  | 3438     | 0.048  | ng/ul# | 100      |
| 26) Benzo(g,h,i)perylene    | 26.79 | 276  | 4416     | 0.071  | ng/ul# | 66       |

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

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Quant Time: Dec 16 02:10:24 2018  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN11918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Sat Dec 15 23:04:58 2018  
Response via : Initial Calibration

