

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061019\  
 Data File : BN006234.D  
 Acq On : 10 Jun 2019 21:45  
 Operator : JU/SJ  
 Sample : K3017-21  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 A51F2

Manual Integrations  
 APPROVED

mohammad  
 6/11/2019 4:28:54 PM

Quant Time: Jun 11 02:17:39 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sun Jun 09 05:10:30 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|-------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.65  | 152  | 4884     | 0.40  | ng/ul  | 0.00     |
| 2) Naphthalene-d8           | 10.42 | 136  | 18080    | 0.40  | ng/ul  | 0.00     |
| 6) Acenaphthene-d10         | 14.30 | 164  | 8978     | 0.40  | ng/ul  | 0.00     |
| 10) Phenanthrene-d10        | 17.05 | 188  | 18105m   | 0.40  | ng/ul  | 0.00     |
| 16) Chrysene-d12            | 21.26 | 240  | 14288m   | 0.40  | ng/ul  | 0.00     |
| 20) Perylene-d12            | 23.49 | 264  | 17989    | 0.40  | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |       |        |          |
| 4) 2-Methylnaphthalene-d10  | 12.02 | 152  | 4890     | 0.18  | ng/ul  | 0.00     |
| 14) Fluoranthene-d10        | 19.09 | 212  | 10464m   | 0.22  | ng/ul  | 0.00     |
| Target Compounds            |       |      |          |       | Ovalue |          |
| 3) Naphthalene              | 10.47 | 128  | 4804     | 0.096 | ng/ul# | 95       |
| 5) 2-Methylnaphthalene      | 12.10 | 142  | 1304     | 0.038 | ng/ul  | 100      |
| 7) Acenaphthylene           | 14.01 | 152  | 18932    | 0.379 | ng/ul  | 99       |
| 8) Acenaphthene             | 14.36 | 153  | 2105     | 0.061 | ng/ul  | 95       |
| 9) Fluorene                 | 15.35 | 166  | 2854     | 0.070 | ng/ul# | 95       |
| 11) Pentachlorophenol       | 16.71 | 266  | 204m     | 0.044 | ng/ul  |          |
| 12) Phenanthrene            | 17.09 | 178  | 257132m  | 4.424 | ng/ul  |          |
| 13) Anthracene              | 17.18 | 178  | 18342m   | 0.330 | ng/ul  |          |
| 15) Fluoranthene            | 19.12 | 202  | 636361m  | 9.916 | ng/ul  |          |
| 17) Pyrene                  | 19.49 | 202  | 428858   | 5.698 | ng/ul  | 99       |
| 18) Benzo(a)anthracene      | 21.24 | 228  | 65154m   | 1.031 | ng/ul  |          |
| 19) Chrysene                | 21.30 | 228  | 233959   | 3.944 | ng/ul  | 99       |
| 21) Benzo(b)fluoranthene    | 22.82 | 252  | 254676m  | 3.395 | ng/ul  |          |
| 22) Benzo(k)fluoranthene    | 22.87 | 252  | 90695m   | 1.241 | ng/ul  |          |
| 23) Benzo(a)pyrene          | 23.39 | 252  | 88205    | 1.263 | ng/ul# | 90       |
| 24) Indeno(1,2,3-cd)pyrene  | 25.72 | 276  | 91980    | 1.216 | ng/ul# | 89       |
| 25) Dibenzo(a,h)anthracene  | 25.73 | 278  | 16978    | 0.281 | ng/ul# | 92       |
| 26) Benzo(g,h,i)perylene    | 26.41 | 276  | 80544    | 1.272 | ng/ul  | 95       |

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

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