

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081820\  
 Data File : BM027130.D  
 Acq On : 18 Aug 2020 17:21  
 Operator : JU/CG  
 Sample : SP5242  
 Misc : SOM02.4 SIM/REG SURROGATE  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SP5242

Manual Integrations  
 APPROVED

mohammad  
 8/19/2020 2:24:44 PM

Quant Time: Aug 18 18:04:26 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM081720.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 18 11:17:38 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|-------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.63  | 152  | 1042     | 0.40  | ng/ul  | 0.00     |
| 2) Naphthalene-d8           | 10.40 | 136  | 3553     | 0.40  | ng/ul  | 0.00     |
| 6) Acenaphthene-d10         | 14.26 | 164  | 2956     | 0.40  | ng/ul  | 0.00     |
| 10) Phenanthrene-d10        | 17.01 | 188  | 5909m    | 0.40  | ng/ul  | 0.00     |
| 16) Chrysene-d12            | 21.21 | 240  | 5230     | 0.40  | ng/ul  | 0.00     |
| 20) Perylene-d12            | 23.39 | 264  | 5971     | 0.40  | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |       |        |          |
| 4) 2-Methylnaphthalene-d10  | 12.00 | 152  | 4779     | 0.80  | ng/ul  | 0.00     |
| 14) Fluoranthene-d10        | 19.04 | 212  | 11467    | 0.68  | ng/ul  | 0.00     |
| Target Compounds            |       |      |          |       | Ovalue |          |
| 3) Naphthalene              | 10.44 | 128  | 13       | 0.001 | ng/ul# | 1        |
| 5) 2-Methylnaphthalene      | 12.14 | 142  | 4        | 0.001 | ng/ul  | 91       |
| 7) Acenaphthylene           | 14.00 | 152  | 3        | 0.000 | ng/ul# | 1        |
| 8) Acenaphthene             | 14.32 | 153  | 7        | 0.001 | ng/ul# | 65       |
| 11) Pentachlorophenol       | 16.67 | 266  | 2        | 0.001 | ng/ul# | 1        |
| 12) Phenanthrene            | 17.05 | 178  | 30       | 0.002 | ng/ul# | 1        |
| 15) Fluoranthene            | 19.07 | 202  | 28       | 0.001 | ng/ul# | 1        |
| 18) Benzo(a)anthracene      | 21.20 | 228  | 77       | 0.004 | ng/ul# | 9        |
| 19) Chrysene                | 21.25 | 228  | 111      | 0.005 | ng/ul# | 34       |
| 21) Benzo(b)fluoranthene    | 22.74 | 252  | 25       | 0.001 | ng/ul# | 1        |
| 22) Benzo(k)fluoranthene    | 22.77 | 252  | 8        | 0.000 | ng/ul# | 1        |
| 24) Indeno(1,2,3-cd)pyrene  | 25.58 | 276  | 5        | 0.000 | ng/ul# | 1        |
| 25) Dibenzo(a,h)anthracene  | 25.60 | 278  | 5        | 0.000 | ng/ul# | 1        |
| 26) Benzo(g,h,i)perylene    | 26.23 | 276  | 5        | 0.000 | ng/ul# | 1        |

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

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