

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM122724\  
 Data File : BM049283.D  
 Acq On : 28 Dec 2024 09:09  
 Operator : RC/JU  
 Sample : P5351-01  
 Misc :  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 E2922

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 12/30/2024  
 Supervised By :mohammad ahmed 12/30/2024

Quant Time: Dec 29 22:13:12 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM121824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Dec 26 11:35:38 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.539  | 152  | 6208     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.298 | 136  | 20035    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.178 | 164  | 12034    | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 16.926 | 188  | 28110m   | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.128 | 240  | 23561m   | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.279 | 264  | 23388m   | 0.400 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.138  | 96   | 20071    | 2.803 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.899 | 152  | 5529     | 0.205 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 18.960 | 212  | 12853m   | 0.192 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       | Qvalue |          |
| 5) Naphthalene              | 10.348 | 128  | 2346     | 0.047 | ng/ul# | 86       |
| 7) 2-Methylnaphthalene      | 11.976 | 142  | 1429     | 0.045 | ng/ul  | 98       |
| 8) 1-Methylnaphthalene      | 12.196 | 142  | 1465     | 0.045 | ng/ul  | 96       |
| 10) Acenaphthylene          | 13.891 | 152  | 2811     | 0.056 | ng/ul# | 87       |
| 11) Acenaphthene            | 14.238 | 153  | 3582     | 0.098 | ng/ul  | 98       |
| 12) Fluorene                | 15.233 | 166  | 6002     | 0.146 | ng/ul  | 97       |
| 15) Phenanthrene            | 16.968 | 178  | 167427m  | 2.280 | ng/ul  |          |
| 16) Anthracene              | 17.057 | 178  | 49063m   | 0.728 | ng/ul  |          |
| 19) Fluoranthene            | 18.993 | 202  | 379843m  | 4.459 | ng/ul  |          |
| 20) Pyrene                  | 19.355 | 202  | 282942m  | 3.114 | ng/ul  |          |
| 21) Benzo(a)anthracene      | 21.113 | 228  | 164481m  | 2.007 | ng/ul  |          |
| 22) Chrysene                | 21.166 | 228  | 151175m  | 1.719 | ng/ul  |          |
| 24) Benzo(b)fluoranthene    | 22.642 | 252  | 167942m  | 2.248 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 22.680 | 252  | 61993m   | 0.756 | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.185 | 252  | 111580m  | 1.557 | ng/ul  |          |
| 27) Indeno(1,2,3-cd)pyrene  | 25.410 | 276  | 77618m   | 0.765 | ng/ul  |          |
| 28) Dibenzo(a,h)anthracene  | 25.413 | 278  | 20630m   | 0.262 | ng/ul  |          |
| 29) Benzo(g,h,i)perylene    | 26.054 | 276  | 76346m   | 0.936 | ng/ul  |          |
| -----                       |        |      |          |       |        |          |

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

