

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN013125\  
 Data File : BN036192.D  
 Acq On : 01 Feb 2025 06:16  
 Operator : RC/JU  
 Sample : Q1224-02  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 A6302

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 02/03/2025  
 Supervised By :mohammad ahmed 02/04/2025

Quant Time: Feb 03 12:01:45 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-SIM-BN012125.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jan 21 23:52:22 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.791  | 152  | 2710     | 0.400 | ng/ul  | -0.03    |
| 4) Naphthalene-d8           | 10.583 | 136  | 7265     | 0.400 | ng/ul  | -0.02    |
| 9) Acenaphthene-d10         | 14.431 | 164  | 4401     | 0.400 | ng/ul  | -0.02    |
| 13) Phenanthrene-d10        | 17.178 | 188  | 9660     | 0.400 | ng/ul  | -0.01    |
| 17) Chrysene-d12            | 21.366 | 240  | 10363    | 0.400 | ng/ul  | # 0.00   |
| 23) Perylene-d12            | 23.657 | 264  | 10603    | 0.400 | ng/ul  | # 0.00   |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.255  | 96   | 7219     | 2.433 | ng/ul  | -0.02    |
| 6) 2-Methylnaphthalene-d10  | 12.178 | 152  | 1073     | 0.117 | ng/ul  | -0.02    |
| 18) Fluoranthene-d10        | 19.210 | 212  | 1763     | 0.062 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 12) Fluorene                | 15.482 | 166  | 387      | 0.023 | ng/ul# | 91       |
| 15) Phenanthrene            | 17.221 | 178  | 22678    | 0.802 | ng/ul  | 97       |
| 16) Anthracene              | 17.309 | 178  | 1861     | 0.071 | ng/ul# | 91       |
| 19) Fluoranthene            | 19.238 | 202  | 96858    | 2.459 | ng/ul  | 100      |
| 20) Pyrene                  | 19.601 | 202  | 79323    | 1.929 | ng/ul  | 100      |
| 21) Benzo(a)anthracene      | 21.348 | 228  | 32928    | 0.890 | ng/ul  | 99       |
| 22) Chrysene                | 21.401 | 228  | 49201    | 1.302 | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 22.967 | 252  | 70171m   | 1.983 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 23.008 | 252  | 26188m   | 0.704 | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.558 | 252  | 41197    | 1.289 | ng/ul  | 97       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.988 | 276  | 30430    | 0.720 | ng/ul# | 95       |
| 28) Dibenzo(a,h)anthracene  | 26.001 | 278  | 6520     | 0.200 | ng/ul# | 82       |
| 29) Benzo(g,h,i)perylene    | 26.712 | 276  | 29284    | 0.892 | ng/ul  | 98       |
| -----                       |        |      |          |       |        |          |

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

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