

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060925\  
 Data File : BP024886.D  
 Acq On : 09 Jun 2025 20:59  
 Operator : RC/JU  
 Sample : Q2260-01  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TP10-MHG-WC

Manual Integrations  
 APPROVED

Reviewed By :Rahul Chavli 06/10/2025  
 Supervised By :Jagrut Upadhyay 06/10/2025

Quant Time: Jun 10 01:32:24 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP060625.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jun 06 16:20:27 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.607  | 152  | 403126   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.378 | 136  | 1646656  | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.242 | 164  | 1022823  | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.042 | 188  | 1950083  | 20.000 | ng    | -0.02    |
| 76) Chrysene-d12            | 21.471 | 240  | 1509343  | 20.000 | ng    | -0.01    |
| 86) Perylene-d12            | 24.724 | 264  | 1704631  | 20.000 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.243  | 112  | 1970441  | 81.589 | ng    | 0.00     |
| 7) Phenol-d6                | 6.813  | 99   | 2515256  | 78.715 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.760  | 82   | 1571895  | 46.387 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.766 | 330  | 1180279  | 83.465 | ng    | -0.02    |
| 45) 2-Fluorobiphenyl        | 12.854 | 172  | 3640737  | 47.951 | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.772 | 244  | 5058513  | 60.067 | ng    | -0.02    |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 17.083 | 178  | 467072   | 4.334  | ng    | 99       |
| 75) Fluoranthene            | 19.177 | 202  | 921237   | 7.376  | ng    | 99       |
| 78) Pyrene                  | 19.554 | 202  | 970321   | 10.290 | ng    | 99       |
| 81) Benzo(a)anthracene      | 21.454 | 228  | 355247   | 3.680  | ng    | 98       |
| 83) Chrysene                | 21.518 | 228  | 317703m  | 3.473  | ng    |          |
| 88) Benzo(b)fluoranthene    | 23.695 | 252  | 383801   | 3.934  | ng    | # 92     |
| 90) Benzo(a)pyrene          | 24.571 | 252  | 323864   | 3.399  | ng    | # 93     |
| 92) Benzo(g,h,i)perylene    | 29.553 | 276  | 207061   | 2.061  | ng    | # 91     |
| -----                       |        |      |          |        |       |          |

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

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