

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121025\  
 Data File : BP026283.D  
 Acq On : 10 Dec 2025 19:56  
 Operator : RC/JU  
 Sample : Q3826-01  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TP-13

Quant Time: Dec 11 03:07:55 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP111825.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Dec 05 14:52:44 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.649  | 152  | 175508   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.413 | 136  | 668328   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.278 | 164  | 421570   | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10        | 17.095 | 188  | 857071   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 21.542 | 240  | 1011033  | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 24.842 | 264  | 1205849  | 20.000 | ng    | 0.02     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.272  | 112  | 668227   | 61.211 | ng    | 0.00     |
| 7) Phenol-d6                | 6.831  | 99   | 808989   | 58.707 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.784  | 82   | 478483   | 40.752 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.807 | 330  | 368546   | 67.224 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.889 | 172  | 1142367  | 39.397 | ng    | -0.01    |
| 79) Terphenyl-d14           | 19.836 | 244  | 1994393  | 40.319 | ng    | 0.00     |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 17.136 | 178  | 189876   | 3.930  | ng    | 99       |
| 75) Fluoranthene            | 19.236 | 202  | 323958   | 5.509  | ng    | 99       |
| 78) Pyrene                  | 19.613 | 202  | 299612   | 4.548  | ng    | 100      |
| 81) Benzo(a)anthracene      | 21.524 | 228  | 152665   | 2.201  | ng    | 98       |
| 83) Chrysene                | 21.589 | 228  | 143701   | 2.188  | ng    | 98       |
| 88) Benzo(b)fluoranthene    | 23.795 | 252  | 190493   | 2.599  | ng    | 99       |
| 90) Benzo(a)pyrene          | 24.695 | 252  | 141159   | 2.031  | ng    | 98       |
| -----                       |        |      |          |        |       |          |

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

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