

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103025\  
 Data File : BP026051.D  
 Acq On : 30 Oct 2025 21:11  
 Operator : RC/JU  
 Sample : Q3466-07  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 B7-1.0-2.0-20251023

Quant Time: Oct 31 02:17:22 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP102925.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Oct 30 11:51:34 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.684  | 152  | 374852   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.454 | 136  | 1574285  | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.319 | 164  | 1003733  | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.137 | 188  | 2090155  | 20.000 | ng    | 0.01     |
| 76) Chrysene-d12              | 21.583 | 240  | 1916756  | 20.000 | ng    | 0.02     |
| 86) Perylene-d12              | 24.918 | 264  | 1966081  | 20.000 | ng    | 0.04     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.302  | 112  | 1428503  | 62.883 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.849  | 99   | 1860557  | 64.347 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 8.813  | 82   | 981703   | 38.874 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol      | 15.837 | 330  | 655937   | 60.448 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 12.931 | 172  | 2308607  | 33.050 | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.878 | 244  | 3754937  | 39.801 | ng    | 0.01     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 75) Fluoranthene              | 19.272 | 202  | 341015   | 2.403  | ng    | 99       |
| 78) Pyrene                    | 19.648 | 202  | 440794   | 3.400  | ng    | 98       |
| 81) Benzo(a)anthracene        | 21.566 | 228  | 333231   | 2.530  | ng    | 95       |
| 83) Chrysene                  | 21.636 | 228  | 350457   | 2.809  | ng    | 98       |
| 84) Bis(2-ethylhexyl)phtha... | 21.542 | 149  | 58759    | 3.317  | ng    | # 93     |
| 88) Benzo(b)fluoranthene      | 23.871 | 252  | 511476   | 4.191  | ng    | # 95     |
| 90) Benzo(a)pyrene            | 24.760 | 252  | 396184   | 3.533  | ng    | # 95     |
| 92) Benzo(g,h,i)perylene      | 29.853 | 276  | 301510   | 2.653  | ng    | 96       |
| -----                         |        |      |          |        |       |          |

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

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