

Data Path : C:\Users\eUser\Desktop\Dhruv Practice\BP\  
 Data File : BP024968.D  
 Acq On : 16 Jun 2025 19:46  
 Operator : RC/JU  
 Sample : Q2324-01 2X  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 HD-01-6132025

Quant Time: Jun 17 01:21:09 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   | 7.613  | 152  | 326630   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.384 | 136  | 1296419  | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.254 | 164  | 802382   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.060 | 188  | 1535221  | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 21.495 | 240  | 1553019  | 20.000 | ng    | 0.01     |
| 86) Perylene-d12            | 24.760 | 264  | 1838346  | 20.000 | ng    | 0.04     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.243  | 112  | 779233   | 39.822 | ng    | 0.00     |
| 7) Phenol-d6                | 6.825  | 99   | 1002613  | 38.725 | ng    | 0.01     |
| 23) Nitrobenzene-d5         | 8.760  | 82   | 610126   | 22.869 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.784 | 330  | 441727   | 39.819 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.860 | 172  | 1441261  | 24.197 | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.783 | 244  | 2146368  | 24.770 | ng    | -0.01    |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene            | 17.107 | 178  | 965453   | 11.380 | ng    | 99       |
| 72) Anthracene              | 17.195 | 178  | 333820   | 3.885  | ng    | 100      |
| 75) Fluoranthene            | 19.195 | 202  | 1767655  | 17.976 | ng    | 100      |
| 78) Pyrene                  | 19.572 | 202  | 1577485  | 16.259 | ng    | 99       |
| 81) Benzo(a)anthracene      | 21.472 | 228  | 849447   | 8.552  | ng    | 99       |
| 83) Chrysene                | 21.536 | 228  | 759715   | 8.072  | ng    | 96       |
| 87) Indeno(1,2,3-cd)pyrene  | 28.489 | 276  | 523741   | 3.905  | ng    | 95       |
| 88) Benzo(b)fluoranthene    | 23.736 | 252  | 1105833m | 10.511 | ng    |          |
| 89) Benzo(k)fluoranthene    | 23.789 | 252  | 296346m  | 2.767  | ng    |          |
| 90) Benzo(a)pyrene          | 24.607 | 252  | 810993   | 7.893  | ng    | 94       |
| 92) Benzo(g,h,i)perylene    | 29.659 | 276  | 500114   | 4.617  | ng    | 94       |
| -----                       |        |      |          |        |       |          |

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

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