

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100725\  
 Data File : BN038047.D  
 Acq On : 07 Oct 2025 15:41  
 Operator : RC/JU  
 Sample : Q3248-07  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BP-VPB-184-GW-660-662

Quant Time: Oct 07 17:11:44 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN092325.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 24 11:00:01 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| -----                         |        |      |          |       |       |          |
| Internal Standards            |        |      |          |       |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.804  | 152  | 7914     | 0.400 | ng    | -0.02    |
| 7) Naphthalene-d8             | 10.605 | 136  | 22866    | 0.400 | ng    | -0.02    |
| 13) Acenaphthene-d10          | 14.452 | 164  | 12377    | 0.400 | ng    | -0.02    |
| 19) Phenanthrene-d10          | 17.198 | 188  | 24052    | 0.400 | ng    | #-0.02   |
| 29) Chrysene-d12              | 21.393 | 240  | 13934    | 0.400 | ng    | 0.00     |
| 35) Perylene-d12              | 23.709 | 264  | 10696    | 0.400 | ng    | #-0.02   |
| System Monitoring Compounds   |        |      |          |       |       |          |
| 4) 2-Fluorophenol             | 5.377  | 112  | 2864     | 0.159 | ng    | 0.00     |
| 5) Phenol-d6                  | 6.959  | 99   | 2376     | 0.100 | ng    | -0.01    |
| 8) Nitrobenzene-d5            | 8.950  | 82   | 6210     | 0.356 | ng    | -0.02    |
| 11) 2-Methylnaphthalene-d10   | 12.197 | 152  | 10942    | 0.344 | ng    | -0.02    |
| 14) 2,4,6-Tribromophenol      | 15.945 | 330  | 1637     | 0.349 | ng    | -0.02    |
| 15) 2-Fluorobiphenyl          | 13.073 | 172  | 18912    | 0.373 | ng    | -0.02    |
| 27) Fluoranthene-d10          | 19.239 | 212  | 23036    | 0.381 | ng    | -0.01    |
| 31) Terphenyl-d14             | 19.838 | 244  | 20900    | 0.753 | ng    | -0.01    |
| Target Compounds              |        |      |          |       |       |          |
| 2) 1,4-Dioxane                | 3.290  | 88   | 1097     | 0.137 | ng    | # 60     |
| 33) Chrysene                  | 21.429 | 228  | 1196     | 0.023 | ng    | # 87     |
| 34) Bis(2-ethylhexyl)phtha... | 21.313 | 149  | 2923     | 0.152 | ng    | # 97     |
| 36) Indeno(1,2,3-cd)pyrene    | 26.075 | 276  | 1214     | 0.025 | ng    | # 84     |
| 41) Benzo(g,h,i)perylene      | 26.794 | 276  | 1042     | 0.026 | ng    | # 57     |
| -----                         |        |      |          |       |       |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100725\  
Data File : BN038047.D  
Acq On : 07 Oct 2025 15:41  
Operator : RC/JU  
Sample : Q3248-07  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BP-VPB-184-GW-660-662

Quant Time: Oct 07 17:11:44 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN092325.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Wed Sep 24 11:00:01 2025  
Response via : Initial Calibration

