

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061325\  
 Data File : BP024944.D  
 Acq On : 13 Jun 2025 14:51  
 Operator : RC/JU  
 Sample : Q2286-01 10X  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 LAW-25-0082-LAW-250083

Quant Time: Jun 13 15:17:31 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   | 7.608  | 152  | 245599   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.372 | 136  | 961065   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.254 | 164  | 595173   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.066 | 188  | 1031307  | 20.000 | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.489 | 240  | 1352680  | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 24.742 | 264  | 1586030  | 20.000 | ng    | 0.02     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.249  | 112  | 87793    | 5.967  | ng    | 0.00     |
| 7) Phenol-d6                | 6.837  | 99   | 107544   | 5.524  | ng    | 0.02     |
| 23) Nitrobenzene-d5         | 8.772  | 82   | 75392    | 3.812  | ng    | 0.01     |
| 42) 2,4,6-Tribromophenol    | 15.784 | 330  | 66256    | 8.052  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.854 | 172  | 203709   | 4.611  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.789 | 244  | 337905   | 4.477  | ng    | 0.00     |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 31) Naphthalene             | 10.431 | 128  | 920419   | 18.688 | ng    | 100      |
| 37) 2-Methylnaphthalene     | 12.049 | 142  | 345230   | 11.051 | ng    | 98       |
| 38) 1-Methylnaphthalene     | 12.266 | 142  | 171931   | 5.148  | ng    | 99       |
| 46) 1,1'-Biphenyl           | 13.072 | 154  | 90424    | 2.091  | ng    | 98       |
| 52) Acenaphthene            | 14.319 | 154  | 140835   | 4.434  | ng    | # 89     |
| 55) Dibenzofuran            | 14.660 | 168  | 243994   | 4.786  | ng    | # 54     |
| 58) Fluorene                | 15.325 | 166  | 269532   | 6.544  | ng    | # 79     |
| 71) Phenanthrene            | 17.113 | 178  | 1480436  | 25.976 | ng    | # 85     |
| 75) Fluoranthene            | 19.207 | 202  | 510102   | 7.722  | ng    | 81       |
| 78) Pyrene                  | 19.584 | 202  | 323201   | 3.825  | ng    | # 80     |
| -----                       |        |      |          |        |       |          |

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

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