

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111225\  
 Data File : BP026119.D  
 Acq On : 12 Nov 2025 19:21  
 Operator : RC/JU  
 Sample : Q3597-07  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TP-3

Quant Time: Nov 13 01:12:17 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   | 7.672  | 152  | 284678   | 20.000  | ng    | -0.02    |
| 21) Naphthalene-d8          | 10.437 | 136  | 1115121  | 20.000  | ng    | -0.02    |
| 39) Acenaphthene-d10        | 14.319 | 164  | 695657   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.119 | 188  | 1409983  | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.583 | 240  | 1541508  | 20.000  | ng    | 0.02     |
| 86) Perylene-d12            | 24.912 | 264  | 1739307  | 20.000  | ng    | 0.03     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.296  | 112  | 1592022  | 92.279  | ng    | 0.00     |
| 7) Phenol-d6                | 6.849  | 99   | 1975358  | 89.958  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.813  | 82   | 1136340  | 63.526  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 15.836 | 330  | 775282   | 103.086 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.925 | 172  | 2891059  | 59.717  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.877 | 244  | 4603449  | 60.673  | ng    | 0.01     |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

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

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