

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP050925\  
 Data File : BP024588.D  
 Acq On : 09 May 2025 13:03  
 Operator : RC/JU  
 Sample : Q1968-08  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 MH-M

Quant Time: May 09 13:51:04 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP041425.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri May 09 12:07:46 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.699  | 152  | 141319   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.469 | 136  | 536094   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.328 | 164  | 317044   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.133 | 188  | 623184   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.574 | 240  | 736576   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 24.898 | 264  | 881896   | 20.000  | ng    | 0.01     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.328  | 112  | 1021387  | 119.504 | ng    | 0.00     |
| 7) Phenol-d6                | 6.893  | 99   | 1102643  | 94.218  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.846  | 82   | 857918   | 91.252  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.845 | 330  | 707020   | 161.182 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.940 | 172  | 1881525  | 90.865  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.857 | 244  | 3290893  | 90.055  | ng    | 0.00     |

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

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

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