

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100825\
 Data File : BP025959.D
 Acq On : 08 Oct 2025 18:51
 Operator : RC/JU
 Sample : Q3308-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 S-3(6.0-6.5)

Quant Time: Oct 09 01:28:49 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 08 16:22:13 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.737	152	114937	20.000	ng	0.00
21) Naphthalene-d8	10.507	136	426606	20.000	ng	0.02
39) Acenaphthene-d10	14.348	164	286180	20.000	ng	0.01
64) Phenanthrene-d10	17.148	188	599775	20.000	ng	0.02
76) Chrysene-d12	21.589	240	693716	20.000	ng	0.02
86) Perylene-d12	24.918	264	790565	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.378	112	454839	64.589	ng	0.00
7) Phenol-d6	6.954	99	572578	61.571	ng	0.02
23) Nitrobenzene-d5	8.890	82	329197	34.537	ng	0.01
42) 2,4,6-Tribromophenol	15.877	330	264062	73.886	ng	0.02
45) 2-Fluorobiphenyl	12.960	172	682272	32.026	ng	0.01
79) Terphenyl-d14	19.860	244	1467018	38.132	ng	0.00

Target Compounds	Qvalue

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

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