

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100825\
 Data File : BP025960.D
 Acq On : 08 Oct 2025 19:32
 Operator : RC/JU
 Sample : Q3308-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 S-12(6.0-6.5)

Quant Time: Oct 09 01:29:28 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	7.731	152	128417	20.000	ng	0.00
21) Naphthalene-d8	10.501	136	499788	20.000	ng	0.01
39) Acenaphthene-d10	14.342	164	351286	20.000	ng	0.00
64) Phenanthrene-d10	17.142	188	763845	20.000	ng	0.01
76) Chrysene-d12	21.571	240	824798	20.000	ng	0.00
86) Perylene-d12	24.918	264	868398	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.378	112	408201	51.881	ng	0.00
7) Phenol-d6	6.949	99	523929	50.426	ng	0.01
23) Nitrobenzene-d5	8.884	82	293489	26.282	ng	0.00
42) 2,4,6-Tribromophenol	15.866	330	263323	60.024	ng	0.01
45) 2-Fluorobiphenyl	12.954	172	628312	24.027	ng	0.00
79) Terphenyl-d14	19.848	244	1486774	32.504	ng	0.00

Target Compounds	Qvalue

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

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