

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100925\
 Data File : BP025966.D
 Acq On : 09 Oct 2025 15:46
 Operator : RC/JU
 Sample : Q3308-01RE
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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

Quant Time: Oct 09 16:08:16 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	118110	20.000	ng	0.00
21) Naphthalene-d8	10.501	136	448145	20.000	ng	0.01
39) Acenaphthene-d10	14.336	164	312575	20.000	ng	0.00
64) Phenanthrene-d10	17.136	188	686694	20.000	ng	0.00
76) Chrysene-d12	21.583	240	796537	20.000	ng	0.01
86) Perylene-d12	24.924	264	827201	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.378	112	477340	65.963	ng	0.00
7) Phenol-d6	6.943	99	593476	62.104	ng	0.00
23) Nitrobenzene-d5	8.884	82	338756	33.831	ng	0.00
42) 2,4,6-Tribromophenol	15.854	330	302797	77.570	ng	0.00
45) 2-Fluorobiphenyl	12.954	172	735122	31.593	ng	0.00
79) Terphenyl-d14	19.854	244	1691318	38.287	ng	0.00

Target Compounds	Qvalue

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

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