

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081225\
 Data File : BP025368.D
 Acq On : 12 Aug 2025 15:40
 Operator : CG/JU
 Sample : PB169149TB
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB169149TB

Quant Time: Aug 12 16:02:59 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 06 06:41:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.802	152	216198	20.000	ng	0.00
21) Naphthalene-d8	10.572	136	822002	20.000	ng	0.00
39) Acenaphthene-d10	14.425	164	481171	20.000	ng	0.00
64) Phenanthrene-d10	17.225	188	966595	20.000	ng	0.00
76) Chrysene-d12	21.666	240	1099080	20.000	ng	0.00
86) Perylene-d12	25.059	264	1180748	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.413	112	1602688	119.229	ng	0.00
7) Phenol-d6	6.972	99	1882681	106.421	ng	0.00
23) Nitrobenzene-d5	8.937	82	1217011	82.024	ng	0.00
42) 2,4,6-Tribromophenol	15.936	330	679469	141.549	ng	0.01
45) 2-Fluorobiphenyl	13.037	172	2723558	78.180	ng	0.00
79) Terphenyl-d14	19.954	244	4447639	76.433	ng	0.01

Target Compounds	Qvalue

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

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