

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081225\
 Data File : BP025363.D
 Acq On : 12 Aug 2025 12:12
 Operator : CG/JU
 Sample : PB169180BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB169180BL

Quant Time: Aug 12 12:35:00 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	7.796	152	223850	20.000	ng	0.00
21) Naphthalene-d8	10.572	136	834408	20.000	ng	0.00
39) Acenaphthene-d10	14.419	164	478110	20.000	ng	0.00
64) Phenanthrene-d10	17.225	188	927445	20.000	ng	0.00
76) Chrysene-d12	21.671	240	966084	20.000	ng	0.01
86) Perylene-d12	25.054	264	1157346	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.408	112	1730160	124.312	ng	0.00
7) Phenol-d6	6.972	99	2002834	109.343	ng	0.00
23) Nitrobenzene-d5	8.943	82	1296349	86.072	ng	0.00
42) 2,4,6-Tribromophenol	15.937	330	691825	145.046	ng	0.01
45) 2-Fluorobiphenyl	13.037	172	2924204	84.477	ng	0.00
79) Terphenyl-d14	19.948	244	4130131	80.748	ng	0.00

Target Compounds	Qvalue

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

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