

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080825\
 Data File : BP025329.D
 Acq On : 08 Aug 2025 12:19
 Operator : CG/JU
 Sample : PB169122TB
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB169122TB

Quant Time: Aug 08 12:46:10 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	268110	20.000	ng	0.00
21) Naphthalene-d8	10.578	136	1066396	20.000	ng	0.00
39) Acenaphthene-d10	14.419	164	651726	20.000	ng	0.00
64) Phenanthrene-d10	17.213	188	1241724	20.000	ng	-0.01
76) Chrysene-d12	21.672	240	1038131	20.000	ng	0.01
86) Perylene-d12	25.065	264	1111724	20.000	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.408	112	1998847	119.909	ng	0.00
7) Phenol-d6	6.966	99	2493898	113.676	ng	0.00
23) Nitrobenzene-d5	8.943	82	1555744	80.824	ng	0.00
42) 2,4,6-Tribromophenol	15.931	330	866192	133.225	ng	0.00
45) 2-Fluorobiphenyl	13.037	172	3529574	74.803	ng	0.00
79) Terphenyl-d14	19.948	244	4531870	82.453	ng	0.00

Target Compounds	Qvalue

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

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