

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061625\
 Data File : BP024958.D
 Acq On : 16 Jun 2025 12:55
 Operator : RC/JU
 Sample : PB168456BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB168456BL

Quant Time: Jun 16 13:50:26 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 06 16:20:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	325815	20.000	ng	0.00
21) Naphthalene-d8	10.378	136	1221526	20.000	ng	0.00
39) Acenaphthene-d10	14.254	164	737769	20.000	ng	0.00
64) Phenanthrene-d10	17.054	188	1294055	20.000	ng	0.00
76) Chrysene-d12	21.489	240	1084530	20.000	ng	0.00
86) Perylene-d12	24.747	264	1309880	20.000	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.243	112	2881068	147.602	ng	0.00
7) Phenol-d6	6.819	99	3475271	134.565	ng	0.00
23) Nitrobenzene-d5	8.760	82	2160111	85.931	ng	0.00
42) 2,4,6-Tribromophenol	15.777	330	1356504	132.990	ng	0.00
45) 2-Fluorobiphenyl	12.854	172	4535780	82.821	ng	0.00
79) Terphenyl-d14	19.783	244	5491241	90.746	ng	-0.01

Target Compounds	Qvalue

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

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