

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061825\
 Data File : BP024988.D
 Acq On : 18 Jun 2025 10:36
 Operator : RC/JU
 Sample : PB168509BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB168509BL

Quant Time: Jun 18 11:35:51 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	7.608	152	278843	20.000	ng	0.00
21) Naphthalene-d8	10.378	136	1046607	20.000	ng	0.00
39) Acenaphthene-d10	14.260	164	632257	20.000	ng	0.01
64) Phenanthrene-d10	17.072	188	1245796	20.000	ng	0.01
76) Chrysene-d12	21.501	240	1328391	20.000	ng	0.02
86) Perylene-d12	24.771	264	1595374	20.000	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.243	112	2391475	143.158	ng	0.00
7) Phenol-d6	6.819	99	2902428	131.316	ng	0.00
23) Nitrobenzene-d5	8.761	82	1812094	84.134	ng	0.00
42) 2,4,6-Tribromophenol	15.790	330	1178919	134.868	ng	0.00
45) 2-Fluorobiphenyl	12.854	172	3825055	81.499	ng	0.00
79) Terphenyl-d14	19.795	244	6242246	84.220	ng	0.00

Target Compounds	Qvalue

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

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