

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090924\
 Data File : BP021868.D
 Acq On : 09 Sep 2024 20:11
 Operator : RC/JU
 Sample : PB163214BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB163214BL

Quant Time: Sep 10 00:31:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP090924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 00:09:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	214440	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	817062	20.000	ng	0.00
39) Acenaphthene-d10	14.387	164	482717	20.000	ng	0.00
64) Phenanthrene-d10	17.175	188	741617	20.000	ng	0.00
76) Chrysene-d12	21.604	240	683062	20.000	ng	-0.01
86) Perylene-d12	24.945	264	771383	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2224418	141.270	ng	0.00
7) Phenol-d6	6.940	99	2972232	138.449	ng	0.00
23) Nitrobenzene-d5	8.905	82	1775479	114.559	ng	0.00
42) 2,4,6-Tribromophenol	15.893	330	614668	108.230	ng	0.00
45) 2-Fluorobiphenyl	12.993	172	3077110	103.887	ng	0.00
79) Terphenyl-d14	19.892	244	3593905	107.985	ng	0.00

Target Compounds	Qvalue

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

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