

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112222\
 Data File : BP012701.D
 Acq On : 22 Nov 2022 14:51
 Operator : CG/JU
 Sample : PB149038TB
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB149038TB

Quant Time: Nov 23 06:24:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 00:38:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	322171	20.000	ng	0.00
21) Naphthalene-d8	10.857	136	1324724	20.000	ng	0.00
39) Acenaphthene-d10	14.675	164	849677	20.000	ng	0.00
64) Phenanthrene-d10	17.434	188	1880853	20.000	ng	0.00
76) Chrysene-d12	21.516	240	1740935	20.000	ng	0.00
86) Perylene-d12	24.045	264	1652354	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2512528	168.955	ng	0.00
7) Phenol-d6	7.187	99	3318696	155.611	ng	0.00
23) Nitrobenzene-d5	9.205	82	2052849	86.424	ng	0.00
42) 2,4,6-Tribromophenol	16.163	330	1375904	125.791	ng	0.00
45) 2-Fluorobiphenyl	13.304	172	5279799	88.527	ng	0.00
79) Terphenyl-d14	19.981	244	8771793	107.088	ng	0.00

Target Compounds	Qvalue

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

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