

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111822\
 Data File : BP012667.D
 Acq On : 18 Nov 2022 09:57
 Operator : CG/JU
 Sample : PB149019BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB149019BL

Quant Time: Nov 18 10:32:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 18 00:16:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.046	152	258781	20.000	ng	0.00
21) Naphthalene-d8	10.863	136	1131626	20.000	ng	0.00
39) Acenaphthene-d10	14.681	164	771456	20.000	ng	0.00
64) Phenanthrene-d10	17.434	188	1788947	20.000	ng	0.00
76) Chrysene-d12	21.516	240	1943967	20.000	ng	0.00
86) Perylene-d12	24.045	264	2016098	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.587	112	2139940	149.446	ng	0.00
7) Phenol-d6	7.187	99	2841757	142.499	ng	0.00
23) Nitrobenzene-d5	9.211	82	1578067	87.680	ng	0.00
42) 2,4,6-Tribromophenol	16.169	330	1223911	138.632	ng	0.00
45) 2-Fluorobiphenyl	13.310	172	4366482	85.164	ng	0.00
79) Terphenyl-d14	19.981	244	8031408	89.265	ng	0.00

Target Compounds	Qvalue

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

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