

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120321\
 Data File : BM033286.D
 Acq On : 03 Dec 2021 14:56
 Operator : CG/JU
 Sample : PB141145BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB141145BL

Quant Time: Dec 03 15:46:30 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 24 15:16:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	123984	20.000	ng	0.00
21) Naphthalene-d8	10.725	136	494362	20.000	ng	#-0.01
39) Acenaphthene-d10	14.554	164	292558	20.000	ng	0.00
64) Phenanthrene-d10	17.289	188	556663	20.000	ng	# 0.00
76) Chrysene-d12	21.448	240	445096	20.000	ng	0.00
86) Perylene-d12	23.777	264	418909	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.519	112	1066058	141.991	ng	0.01
7) Phenol-d6	7.113	99	1381001	137.342	ng	0.02
23) Nitrobenzene-d5	9.090	82	928525	87.194	ng	-0.01
42) 2,4,6-Tribromophenol	16.036	330	343794	107.801	ng	0.00
45) 2-Fluorobiphenyl	13.178	172	1763248	84.999	ng	0.00
79) Terphenyl-d14	19.895	244	2115481	93.983	ng	-0.01

Target Compounds	Qvalue

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

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