

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120321\
 Data File : BM033301.D
 Acq On : 04 Dec 2021 00:01
 Operator : CG/JU
 Sample : M4917-12
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-28

Quant Time: Dec 06 10:48:59 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	7.925	152	138154	20.000	ng	-0.01
21) Naphthalene-d8	10.725	136	536958	20.000	ng	#-0.01
39) Acenaphthene-d10	14.548	164	326078	20.000	ng	-0.01
64) Phenanthrene-d10	17.289	188	663449	20.000	ng	# 0.00
76) Chrysene-d12	21.448	240	658689	20.000	ng	# 0.00
86) Perylene-d12	23.777	264	663372	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.513	112	504806	60.340	ng	0.00
7) Phenol-d6	7.107	99	401021	35.791	ng	0.01
23) Nitrobenzene-d5	9.090	82	1068809	92.405	ng	-0.01
42) 2,4,6-Tribromophenol	16.036	330	433434	121.938	ng	0.00
45) 2-Fluorobiphenyl	13.172	172	1956304	84.611	ng	-0.01
79) Terphenyl-d14	19.895	244	2725353	81.815	ng	-0.01

Target Compounds	Qvalue

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

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