

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
 Data File : BM033309.D
 Acq On : 06 Dec 2021 13:10
 Operator : CG/JU
 Sample : M4917-15
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-31

Quant Time: Dec 06 14:08:33 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 06 13:01:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	116719	20.000	ng	0.00
21) Naphthalene-d8	10.719	136	450700	20.000	ng	# 0.00
39) Acenaphthene-d10	14.548	164	270299	20.000	ng	0.00
64) Phenanthrene-d10	17.289	188	551793	20.000	ng	# 0.00
76) Chrysene-d12	21.447	240	568224	20.000	ng	# 0.00
86) Perylene-d12	23.777	264	622045	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.513	112	452957	64.086	ng	0.00
7) Phenol-d6	7.101	99	363512	38.402	ng	0.00
23) Nitrobenzene-d5	9.089	82	982519	101.202	ng	0.00
42) 2,4,6-Tribromophenol	16.036	330	395913	134.367	ng	0.00
45) 2-Fluorobiphenyl	13.172	172	1702129	88.810	ng	0.00
79) Terphenyl-d14	19.895	244	2235291	77.787	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.772	128	78908	3.319	ng	98
38) 1-Methylnaphthalene	12.595	142	44320	2.807	ng	# 97

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

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