

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121721\
 Data File : BM033498.D
 Acq On : 17 Dec 2021 09:15
 Operator : CG/JU
 Sample : PB141481BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB141409BL

Quant Time: Dec 17 10:53:50 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 17:10:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	31249	20.000	ng	0.00
21) Naphthalene-d8	10.748	136	127253	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	83753	20.000	ng	0.00
64) Phenanthrene-d10	17.330	188	157888	20.000	ng	# 0.00
76) Chrysene-d12	21.512	240	92022	20.000	ng	# 0.00
86) Perylene-d12	23.930	264	68182	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.502	112	282936	156.235	ng	0.00
7) Phenol-d6	7.107	99	363865	149.719	ng	0.00
23) Nitrobenzene-d5	9.107	82	280219	94.551	ng	0.00
42) 2,4,6-Tribromophenol	16.071	330	114282	121.302	ng	0.00
45) 2-Fluorobiphenyl	13.207	172	558233	92.200	ng	0.00
79) Terphenyl-d14	19.954	244	624981	106.923	ng	0.00

Target Compounds	Qvalue

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

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