

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112421\
 Data File : BM033252.D
 Acq On : 24 Nov 2021 18:12
 Operator : CG/JU
 Sample : M4692-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WW-EFF

Quant Time: Nov 26 04:09:38 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	7.937	152	128546	20.000	ng	0.00
21) Naphthalene-d8	10.736	136	514451	20.000	ng	# 0.00
39) Acenaphthene-d10	14.560	164	328303	20.000	ng	0.00
64) Phenanthrene-d10	17.307	188	665873	20.000	ng	# 0.01
76) Chrysene-d12	21.465	240	653114	20.000	ng	0.01
86) Perylene-d12	23.806	264	637503	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.513	112	387463	49.776	ng	0.00
7) Phenol-d6	7.113	99	321640	30.852	ng	0.02
23) Nitrobenzene-d5	9.101	82	1109050	100.079	ng	0.00
42) 2,4,6-Tribromophenol	16.048	330	358148	100.075	ng	0.00
45) 2-Fluorobiphenyl	13.183	172	2109012	90.598	ng	0.00
79) Terphenyl-d14	19.912	244	2144436	64.926	ng	0.00
Target Compounds						
32) Benzoic acid	10.166	122	1225707	252.822	ng	Qvalue 88

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

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