

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012021\
 Data File : BM028477.D
 Acq On : 20 Jan 2021 20:44
 Operator : CG/JU
 Sample : M1141-23RE
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETGI-354RE

Quant Time: Jan 21 00:34:10 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 20 14:45:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	34658	20.00	ng	0.00
21) Naphthalene-d8	10.845	136	132764	20.00	ng	# 0.00
39) Acenaphthene-d10	14.669	164	69483	20.00	ng	0.00
64) Phenanthrene-d10	17.398	188	136234	20.00	ng	# 0.00
76) Chrysene-d12	21.527	240	126868	20.00	ng	# 0.00
86) Perylene-d12	23.815	264	128256	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.557	112	276671	126.10	ng	0.00
7) Phenol-d6	7.193	99	321487	107.62	ng	0.00
23) Nitrobenzene-d5	9.210	82	230619	82.39	ng	0.00
42) 2,4,6-Tribromophenol	16.151	330	123165	133.79	ng	0.00
45) 2-Fluorobiphenyl	13.298	172	466489	84.23	ng	0.00
79) Terphenyl-d14	19.992	244	580033	82.56	ng	0.00

Target Compounds	Qvalue

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

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