

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061021\
 Data File : BM030381.D
 Acq On : 10 Jun 2021 19:15
 Operator : CG/JU
 Sample : M2613-22
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(118-122)

Quant Time: Jun 11 07:53:30 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 20:51:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	75635	20.000	ng	0.00
21) Naphthalene-d8	10.633	136	333802	20.000	ng	# 0.00
39) Acenaphthene-d10	14.468	164	277646	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	793292	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1166853	20.000	ng	0.00
86) Perylene-d12	23.638	264	1213937	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	22619	4.862	ng	0.00
7) Phenol-d6	7.010	99	215916	32.058	ng	0.00
23) Nitrobenzene-d5	8.998	82	224178	32.718	ng	-0.01
42) 2,4,6-Tribromophenol	15.957	330	3375	0.967	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	1466257	69.449	ng	0.00
79) Terphenyl-d14	19.815	244	4764717	71.722	ng	0.00

Target Compounds	Qvalue

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

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