

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
 Data File : BM030407.D
 Acq On : 11 Jun 2021 16:39
 Operator : CG/JU
 Sample : M2613-22
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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

Quant Time: Jun 11 17:10:41 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 14:10:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	165379	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	589234	20.000	ng	# 0.00
39) Acenaphthene-d10	14.457	164	334920	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	650259	20.000	ng	# 0.00
76) Chrysene-d12	21.356	240	664786	20.000	ng	0.00
86) Perylene-d12	23.638	264	784944	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	486539	47.826	ng	0.00
7) Phenol-d6	6.998	99	581286	39.471	ng	0.00
23) Nitrobenzene-d5	8.992	82	939374	77.667	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	427343	101.486	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	2003545	78.669	ng	0.00
79) Terphenyl-d14	19.809	244	2942014	77.731	ng	0.00

Target Compounds	Qvalue

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

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