

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061421\
 Data File : BM030448.D
 Acq On : 14 Jun 2021 22:05
 Operator : CG/JU
 Sample : M2677-17
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PIER1-2

Quant Time: Jun 15 03:51:42 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	7.833	152	168945	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	617655	20.000	ng	# 0.00
39) Acenaphthene-d10	14.457	164	375291	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	787950	20.000	ng	0.00
76) Chrysene-d12	21.350	240	906715	20.000	ng	0.00
86) Perylene-d12	23.627	264	971405	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	928300	89.325	ng	0.01
7) Phenol-d6	7.010	99	1520348	101.057	ng	0.01
23) Nitrobenzene-d5	8.986	82	1040055	82.034	ng	0.00
42) 2,4,6-Tribromophenol	15.939	330	347532	73.654	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	2376693	83.282	ng	0.00
79) Terphenyl-d14	19.803	244	4154399	80.477	ng	0.00

Target Compounds	Qvalue

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

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