

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070121\
 Data File : BF124510.D
 Acq On : 1 Jul 2021 14:22
 Operator : JU/CG
 Sample : M2875-12DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-6DL

Quant Time: Jul 01 14:52:31 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 01 14:15:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.681	152	29852	20.000	ng	0.00
21) Naphthalene-d8	7.963	136	109157	20.000	ng	# 0.00
39) Acenaphthene-d10	9.716	164	67289	20.000	ng	0.00
64) Phenanthrene-d10	11.204	188	119473	20.000	ng	0.00
76) Chrysene-d12	13.851	240	70477	20.000	ng	0.00
86) Perylene-d12	15.280	264	78260	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.316	112	42594	22.440	ng	0.01
7) Phenol-d6	6.345	99	29737	11.998	ng	0.00
23) Nitrobenzene-d5	7.251	82	98731	37.502	ng	0.00
42) 2,4,6-Tribromophenol	10.510	330	44670	55.565	ng	0.00
45) 2-Fluorobiphenyl	9.039	172	200543	36.509	ng	0.00
79) Terphenyl-d14	12.786	244	196496	41.760	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.363	94	10721	4.001	ng	# 1
17) 2-Methylphenol	6.963	107	12206	7.231	ng	# 54
27) 2,4-Dimethylphenol	7.651	122	102619	60.181	ng	86
31) Naphthalene	7.986	128	308265	48.001	ng	97
37) 2-Methylnaphthalene	8.681	142	85313	18.773	ng	97
38) 1-Methylnaphthalene	8.775	142	51169	12.105	ng	92

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

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