

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060122\
 Data File : BF128634.D
 Acq On : 01 Jun 2022 20:52
 Operator : CG\JU
 Sample : N3087-06DL 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-06DL

Quant Time: Jun 02 09:37:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 27 01:51:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	102522	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	313523	20.000	ng	-0.01
39) Acenaphthene-d10	9.869	164	156163	20.000	ng	-0.01
64) Phenanthrene-d10	11.357	188	236323	20.000	ng	#-0.01
76) Chrysene-d12	13.998	240	250908	20.000	ng	#-0.01
86) Perylene-d12	15.462	264	233828	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	129830	20.976	ng	0.05
7) Phenol-d6	6.528	99	96955	12.286	ng	0.06
23) Nitrobenzene-d5	7.398	82	239225	44.017	ng	-0.01
42) 2,4,6-Tribromophenol	10.663	330	96133	64.675	ng	-0.01
45) 2-Fluorobiphenyl	9.192	172	475728	49.492	ng	-0.01
79) Terphenyl-d14	12.939	244	460652	35.470	ng	-0.02
Target Compounds						
						Qvalue
10) Phenol	6.545	94	35554	4.547	ng	84
17) 2-Methylphenol	7.128	107	33885	6.203	ng	# 93
22) Acetophenone	7.239	105	331086	45.798	ng	# 92
27) 2,4-Dimethylphenol	7.792	122	144266	32.321	ng	93
31) Naphthalene	8.139	128	750221	48.890	ng	99
37) 2-Methylnaphthalene	8.828	142	133612	13.795	ng	98
38) 1-Methylnaphthalene	8.928	142	100363	10.654	ng	100

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

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