

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032922\
 Data File : BF127563.D
 Acq On : 29 Mar 2022 14:10
 Operator : CG\JU
 Sample : N2154-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-5

Quant Time: Mar 29 15:56:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF032122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 29 07:13:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	81758	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	301766	20.000	ng	0.00
39) Acenaphthene-d10	9.875	164	149461	20.000	ng	0.00
64) Phenanthrene-d10	11.369	188	216532	20.000	ng	0.00
76) Chrysene-d12	14.016	240	177590	20.000	ng	0.00
86) Perylene-d12	15.504	264	195421	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	230811	45.063	ng	0.00
7) Phenol-d6	6.487	99	181638	25.943	ng	0.00
23) Nitrobenzene-d5	7.404	82	643420	91.142	ng	0.00
42) 2,4,6-Tribromophenol	10.675	330	178517	112.124	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	993641	97.079	ng	0.00
79) Terphenyl-d14	12.951	244	693053	63.885	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.504	94	18311	2.402	ng	82
31) Naphthalene	8.139	128	41460	2.588	ng	# 95
32) Benzoic acid	7.975	122	137058	55.121	ng	94

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

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