

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121522\
 Data File : BF131654.D
 Acq On : 15 Dec 2022 17:52
 Operator : CG\JU
 Sample : N6001-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-7

Quant Time: Dec 16 01:53:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 16 01:44:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	66147	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	239539	20.000	ng	0.00
39) Acenaphthene-d10	9.933	164	139516	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	271131	20.000	ng	0.00
76) Chrysene-d12	14.074	240	156542	20.000	ng	# 0.00
86) Perylene-d12	15.580	264	144217	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	233892	58.990	ng	0.00
7) Phenol-d6	6.504	99	225065	43.317	ng	-0.01
23) Nitrobenzene-d5	7.451	82	553964	87.433	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	216540	139.697	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	952496	86.082	ng	0.00
79) Terphenyl-d14	13.021	244	650209	56.962	ng	0.00
Target Compounds						
10) Phenol	6.516	94	12611	2.176	ng	Qvalue 96
22) Acetophenone	7.287	105	51230	6.825	ng	# 91
27) 2,4-Dimethylphenol	7.816	122	52081	15.592	ng	97

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

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