

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091521\
 Data File : BF125510.D
 Acq On : 15 Sep 2021 18:22
 Operator : JU/CG
 Sample : M3728-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 N42049

Quant Time: Sep 16 03:06:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF091521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 15 15:08:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	173693	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	665095	20.000	ng	# 0.00
39) Acenaphthene-d10	9.910	164	361404	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	626594	20.000	ng	0.00
76) Chrysene-d12	14.045	240	452536	20.000	ng	# 0.00
86) Perylene-d12	15.527	264	347243	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	598247	54.727	ng	0.00
7) Phenol-d6	6.498	99	481602	33.148	ng	-0.01
23) Nitrobenzene-d5	7.433	82	1029970	76.461	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	519960	118.128	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1804807	73.279	ng	0.00
79) Terphenyl-d14	12.992	244	2080682	74.555	ng	0.00
Target Compounds						
22) Acetophenone	7.269	105	86287	5.264	ng	# 94

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

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