

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090921\
 Data File : BF125407.D
 Acq On : 9 Sep 2021 11:08
 Operator : JU/CG
 Sample : M3646-13DL 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1992DL

Quant Time: Sep 09 11:37:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 08 14:14:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.886	152	137098	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	550853	20.000	ng	# 0.00
39) Acenaphthene-d10	9.927	164	279795	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	499030	20.000	ng	0.00
76) Chrysene-d12	14.057	240	394608	20.000	ng	# 0.00
86) Perylene-d12	15.545	264	387114	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	156998	17.333	ng	0.00
7) Phenol-d6	6.516	99	127977	10.925	ng	-0.01
23) Nitrobenzene-d5	7.445	82	379704	34.702	ng	-0.01
42) 2,4,6-Tribromophenol	10.716	330	115781	41.454	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	694450	34.592	ng	0.00
79) Terphenyl-d14	12.998	244	782829	31.604	ng	0.00
Target Compounds						
3) Pyridine	3.469	79	34025	2.884	ng	# 91
20) 3+4-Methylphenols	7.292	107	438647	45.020	ng	# 74

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

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