

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090121\
 Data File : BF125304.D
 Acq On : 1 Sep 2021 13:56
 Operator : JU/CG
 Sample : M3593-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-280

Quant Time: Sep 01 14:36:09 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 24 12:39:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.904	152	173925	20.000	ng	-0.01
21) Naphthalene-d8	8.186	136	686596	20.000	ng	#-0.01
39) Acenaphthene-d10	9.945	164	371271	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	692942	20.000	ng	0.00
76) Chrysene-d12	14.074	240	562427	20.000	ng	# 0.00
86) Perylene-d12	15.574	264	422263	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.539	112	1172950	102.077	ng	0.01
7) Phenol-d6	6.539	99	1368126	92.060	ng	0.00
23) Nitrobenzene-d5	7.469	82	1061845	77.859	ng	-0.01
42) 2,4,6-Tribromophenol	10.739	330	531592	143.437	ng	0.00
45) 2-Fluorobiphenyl	9.269	172	1842080	69.151	ng	0.00
79) Terphenyl-d14	13.027	244	2415072	68.408	ng	0.00
Target Compounds						
58) Fluorene	10.492	166	55622	2.379	ng	# 93
71) Phenanthrene	11.457	178	325122	8.515	ng	98

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

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