

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092321\
 Data File : BF125610.D
 Acq On : 24 Sep 2021 1:01
 Operator : JU/CG
 Sample : M3822-11
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CURB-SIDE

Quant Time: Sep 24 04:20:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 21 19:21:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	230764	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	942512	20.000	ng	# 0.00
39) Acenaphthene-d10	9.933	164	508766	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	876798	20.000	ng	# 0.00
76) Chrysene-d12	14.057	240	637268	20.000	ng	0.00
86) Perylene-d12	15.539	264	542862	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1357668	93.037	ng	0.00
7) Phenol-d6	6.516	99	1685985	89.609	ng	0.00
23) Nitrobenzene-d5	7.451	82	1036356	70.133	ng	-0.01
42) 2,4,6-Tribromophenol	10.716	330	377503	75.755	ng	-0.01
45) 2-Fluorobiphenyl	9.251	172	2115585	73.790	ng	-0.01
79) Terphenyl-d14	13.004	244	2207073	55.927	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	11.439	178	112373	2.282	ng	98
75) Fluoranthene	12.627	202	150677	3.118	ng	99

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

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