

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093021\
 Data File : BF125733.D
 Acq On : 30 Sep 2021 15:03
 Operator : JU/CG
 Sample : M3971-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RP-COMP-3

Quant Time: Sep 30 15:27:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 15:38:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	205227	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	829131	20.000	ng	# 0.00
39) Acenaphthene-d10	9.922	164	437895	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	761165	20.000	ng	# 0.00
76) Chrysene-d12	14.039	240	564908	20.000	ng	0.00
86) Perylene-d12	15.515	264	435192	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1411079	113.282	ng	0.00
7) Phenol-d6	6.510	99	1565929	93.507	ng	0.00
23) Nitrobenzene-d5	7.445	82	1100401	92.571	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	569729	139.544	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	2157738	89.200	ng	0.00
79) Terphenyl-d14	12.992	244	2409624	73.869	ng	0.00
Target Compounds						
77) Benzidine	12.739	184	1431818	96.482	ng	Qvalue 97

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

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