

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092921\
 Data File : BF125719.D
 Acq On : 29 Sep 2021 20:55
 Operator : JU/CG
 Sample : M3964-17
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CC-17

Quant Time: Sep 30 03:28:43 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	6.886	152	256252	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	995840	20.000	ng	# 0.00
39) Acenaphthene-d10	9.921	164	481172	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	697260	20.000	ng	# 0.00
76) Chrysene-d12	14.039	240	527615	20.000	ng	0.00
86) Perylene-d12	15.515	264	484155	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1493548	96.027	ng	0.00
7) Phenol-d6	6.510	99	1843213	88.149	ng	0.00
23) Nitrobenzene-d5	7.445	82	1095438	76.726	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	427512	95.293	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	2029123	74.049	ng	0.00
79) Terphenyl-d14	12.992	244	1703456	55.912	ng	0.00
Target Compounds						
						Qvalue
32) Benzoic acid	7.857	122	2155	8.797	ng	90
74) Di-n-butylphthalate	11.963	149	174227	4.545	ng	99
80) Butylbenzylphthalate	13.462	149	122974	7.788	ng	97
84) Bis(2-ethylhexyl)phtha...	14.027	149	522725	28.918	ng	100

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

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