

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091019\
 Data File : BF116923.D
 Acq On : 10 Sep 2019 16:06
 Operator : HP/JU
 Sample : PB122853BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122853BL

Quant Time: Sep 10 18:01:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 17:49:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	4858	20.00	ng	0.00
21) Naphthalene-d8	0.00	136	0	0.00	ng	-8.12
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-9.87
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-11.35
76) Chrysene-d12	0.00	240	0	0.00	ng	-13.99
87) Perylene-d12	0.00	264	0	0.00	ng	-15.43
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	669854	2233.84	ng	0.01
7) Phenol-d6	6.47	99	898249	2281.23	ng	0.00
23) Nitrobenzene-d5	7.39	82	575845	0.00	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	198008	0.00	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1034254	0.00	ng	0.00
79) Terphenyl-d14	12.93	244	1018689	0.00	ng	0.00
Target Compounds						
9) Benzaldehyde	6.47	77	1206	4.649	ng	# 2
10) Phenol	6.60	94	1357	3.112	ng	# 1
11) bis(2-Chloroethyl)ether	6.60	93	1027	3.025	ng	# 1
15) Benzyl Alcohol	6.99	79	9626	30.116	ng	# 51
19) n-Nitroso-di-n-propylamine	7.39	70	75186	289.999	ng	# 75

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

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