

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100118\
 Data File : BF109481.D
 Acq On : 1 Oct 2018 19:06
 Operator : JU/SJ
 Sample : J5155-01 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW30S-GW

Quant Time: Oct 02 12:21:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	102208	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	340948	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	163641	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	274762	20.00	ng	0.00
75) Chrysene-d12	13.84	240	245474	20.00	ng	0.00
86) Perylene-d12	15.24	264	185267	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	8546	1.38	ng	0.01
7) Phenol-d6	6.37	99	12431	1.57	ng	0.01
23) Nitrobenzene-d5	7.26	82	35905	6.22	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	9688	6.01	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	68397	6.30	ng	-0.02
78) Terphenyl-d14	12.79	244	41654	4.16	ng	-0.01
Target Compounds						
6) Aniline	6.37	93	30144	2.976	ng	93
20) 3+4-Methylphenols	7.13	107	52264	7.753	ng	90
27) 2,4-Dimethylphenol	7.66	122	66182	14.604	ng	97
31) Naphthalene	8.04	128	7141841	448.259	ng	# 94
37) 2-Methylnaphthalene	8.70	142	519236	50.554	ng	98
45) 1,1'-Biphenyl	9.16	154	75092	5.632	ng	96
51) Acenaphthene	9.77	154	89724	9.236	ng	97
54) Dibenzofuran	9.94	168	110136	7.623	ng	98
57) Fluorene	10.28	166	48040	4.660	ng	99
70) Phenanthrene	11.24	178	59596	4.344	ng	97
72) Carbazole	11.45	167	116391	8.407	ng	97

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

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