

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102218\
 Data File : BF110060.D
 Acq On : 22 Oct 2018 20:20
 Operator : JU/SJ
 Sample : J5604-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Oct 23 04:27:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 14:58:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	186421	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	694321	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	269485	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	400931	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	302242	20.00	ng	-0.01
87) Perylene-d12	15.67	264	235265	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1063678	90.44	ng	0.00
7) Phenol-d6	6.59	99	1324651	90.71	ng	-0.01
23) Nitrobenzene-d5	7.53	82	703674	68.27	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	200552	85.96	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	921328	54.55	ng	-0.01
79) Terphenyl-d14	13.09	244	607840	42.23	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	120386	7.635	ng	# 65
20) 3+4-Methylphenols	7.36	107	42195	3.130	ng	# 74
50) Dimethylphthalate	9.72	163	166140	8.707	ng	99
75) Fluoranthene	12.72	202	59314m	2.861	ng	
78) Pyrene	12.95	202	81035	3.650	ng	# 83
88) Benzo(b)fluoranthene	15.22	252	37119m	2.737	ng	
90) Benzo(a)pyrene	15.60	252	30234	2.424	ng	# 84

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

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