

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111536.D
 Acq On : 17 Dec 2018 13:44
 Operator : JU/SJ
 Sample : J6386-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS006-0612-01

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:47:47 PM

Quant Time: Dec 18 04:42:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	145610	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	529102	20.00	ng	-0.01
39) Acenaphthene-d10	9.82	164	220797	20.00	ng	-0.01
64) Phenanthrene-d10	11.30	188	357880	20.00	ng	-0.01
76) Chrysene-d12	13.95	240	490258	20.00	ng	0.00
87) Perylene-d12	15.39	264	421128	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	918244	104.94	ng	0.00
7) Phenol-d6	6.44	99	1112752	95.60	ng	0.00
23) Nitrobenzene-d5	7.35	82	711962	80.13	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	183830	92.94	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1160739	81.20	ng	-0.01
79) Terphenyl-d14	12.90	244	1222489	58.55	ng	0.00
Target Compounds						
10) Phenol	6.45	94	80609	6.217	ng	95
49) Acenaphthylene	9.68	152	46887	2.221	ng	95
50) Dimethylphthalate	9.54	163	105763	6.651	ng	95
71) Phenanthrene	11.33	178	118922	6.449	ng	93
75) Fluoranthene	12.52	202	116871	6.478	ng	96
78) Pyrene	12.74	202	205623	5.949	ng	96
81) Benzo(a)anthracene	13.94	228	71754m	2.692	ng	
83) Chrysene	13.97	228	79603	2.835	ng	95
88) Benzo(b)fluoranthene	14.97	252	68375m	2.785	ng	
90) Benzo(a)pyrene	15.33	252	70513	3.185	ng	# 83
92) Benzo(g,h,i)perylene	17.23	276	41095	2.397	ng	98

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

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