

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111538.D
 Acq On : 17 Dec 2018 14:39
 Operator : JU/SJ
 Sample : J6386-15
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS007-0006-01

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:47:50 PM

Quant Time: Dec 18 04:53:26 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	176189	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	673526	20.00	ng	-0.01
39) Acenaphthene-d10	9.82	164	285771	20.00	ng	-0.01
64) Phenanthrene-d10	11.31	188	430448	20.00	ng	0.00
76) Chrysene-d12	13.94	240	465065	20.00	ng	0.00
87) Perylene-d12	15.39	264	384822	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	915299	86.45	ng	0.00
7) Phenol-d6	6.44	99	1224146	86.92	ng	0.00
23) Nitrobenzene-d5	7.35	82	764180	67.56	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	195027	76.19	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1110751	60.04	ng	-0.01
79) Terphenyl-d14	12.89	244	934489	47.18	ng	0.00
Target Compounds						
10) Phenol	6.45	94	107532	6.854	ng	92
20) 3+4-Methylphenols	7.20	107	27701	2.169	ng	# 58
49) Acenaphthylene	9.69	152	58044	2.125	ng	94
50) Dimethylphthalate	9.54	163	78623	3.820	ng	98
71) Phenanthrene	11.33	178	222619	10.037	ng	94
75) Fluoranthene	12.52	202	187094	8.622	ng	96
78) Pyrene	12.75	202	314373	9.588	ng	92
81) Benzo(a)anthracene	13.93	228	110983m	4.389	ng	
83) Chrysene	13.97	228	114833	4.311	ng	96
88) Benzo(b)fluoranthene	14.97	252	92786m	4.136	ng	
90) Benzo(a)pyrene	15.32	252	85706	4.236	ng	97
92) Benzo(a,h,i)perylene	17.23	276	45086	2.878	ng	96

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

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