

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121218\
 Data File : BF111338.D
 Acq On : 12 Dec 2018 22:56
 Operator : JU/SJ
 Sample : J6189-01 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-05-121118-A

Manual Integrations
 APPROVED

Sohil
 12/13/2018 6:06:48 PM

Quant Time: Dec 13 04:22:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 13:16:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	262733	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	932283	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	366959	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	602643	20.00	ng	0.00
76) Chrysene-d12	13.95	240	412868	20.00	ng	0.00
87) Perylene-d12	15.39	264	324274	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	947766	61.00	ng	0.00
7) Phenol-d6	6.43	99	1171345	55.92	ng	0.00
23) Nitrobenzene-d5	7.35	82	655635	39.64	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	164191	50.76	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1076520	47.10	ng	0.00
79) Terphenyl-d14	12.90	244	829846	44.24	ng	0.00
Target Compounds						
10) Phenol	6.43	94	88425	3.726	ng	93
50) Dimethylphthalate	9.55	163	79505	3.013	ng	# 98
75) Fluoranthene	12.53	202	128348	4.221	ng	98
78) Pyrene	12.75	202	127892	4.014	ng	96
81) Benzo(a)anthracene	13.94	228	80887m	3.502	ng	
83) Chrysene	13.97	228	78507	3.331	ng	98
88) Benzo(b)fluoranthene	14.97	252	79218m	4.296	ng	
90) Benzo(a)pyrene	15.33	252	58886	3.430	ng	# 80
92) Benzo(g,h,i)perylene	17.22	276	32251	2.258	ng	95

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

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