

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
 Data File : BF112084.D
 Acq On : 17 Jan 2019 12:56
 Operator : JU/SJ
 Sample : K1012-11 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-011518-C

Manual Integrations
 APPROVED

Sohil
 1/17/2019 2:55:44 PM

Quant Time: Jan 17 14:15:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 16 18:22:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	283080	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1060356	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	472191	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	740863	20.00	ng	0.00
76) Chrysene-d12	14.03	240	651276	20.00	ng	0.00
87) Perylene-d12	15.50	264	548591	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	179857	11.53	ng	0.00
7) Phenol-d6	6.50	99	237601	12.18	ng	-0.01
23) Nitrobenzene-d5	7.42	82	121221	7.90	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	46420	9.11	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	295654	9.18	ng	0.00
79) Terphenyl-d14	12.96	244	243687	7.96	ng	-0.01
Target Compounds						
37) 2-Methylnaphthalene	8.86	142	93212	2.846	ng	98
38) 1-Methylnaphthalene	8.96	142	65722m	2.088	ng	
71) Phenanthrene	11.41	178	274789	7.370	ng	97
75) Fluoranthene	12.60	202	360257	9.151	ng	99
78) Pyrene	12.83	202	324363	7.549	ng	98
81) Benzo(a)anthracene	14.02	228	179635	4.811	ng	99
83) Chrysene	14.05	228	147445	4.173	ng	97
86) Indeno(1,2,3-cd)pyrene	16.97	276	65995	2.024	ng	95
88) Benzo(b)fluoranthene	15.07	252	177599m	5.703	ng	
89) Benzo(k)fluoranthene	15.09	252	63617m	2.111	ng	
90) Benzo(a)pyrene	15.44	252	126825	4.441	ng	# 94
92) Benzo(a,h,i)perylene	17.43	276	60557	2.423	ng	98

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

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