

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
 Data File : BF112785.D
 Acq On : 1 Mar 2019 2:16
 Operator : JU/SJ
 Sample : K1650-06 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-3(2.5-3.5)

Manual Integrations
 APPROVED

Sohil
 3/1/2019 3:00:37 PM

Quant Time: Mar 01 08:31:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 27 16:08:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	178901	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	653144	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	257318	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	460967	20.00	ng	0.00
76) Chrysene-d12	13.87	240	498359	20.00	ng	0.00
87) Perylene-d12	15.29	264	469163	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	250737	21.80	ng	0.00
7) Phenol-d6	6.36	99	331574	22.95	ng	0.00
23) Nitrobenzene-d5	7.30	82	181617	16.64	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	65789	24.08	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	301777	8.22	ng	0.00
79) Terphenyl-d14	12.83	244	293830	11.43	ng	0.00
Target Compounds						
71) Phenanthrene	11.27	178	139529	6.084	ng	97
75) Fluoranthene	12.45	202	210608	8.571	ng	99
78) Pyrene	12.68	202	219389	5.222	ng	98
81) Benzo(a)anthracene	13.86	228	115092	3.332	ng	98
83) Chrysene	13.90	228	105553	3.120	ng	97
88) Benzo(b)fluoranthene	14.88	252	116685m	3.845	ng	
90) Benzo(a)pyrene	15.22	252	82954	3.090	ng	98

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

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