

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
 Data File : BF112813.D
 Acq On : 1 Mar 2019 21:20
 Operator : JU/SJ
 Sample : K1675-17 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Z-4-A-B

Manual Integrations
 APPROVED

Sohil
 3/4/2019 4:55:38 PM

Quant Time: Mar 02 00:37:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	157278	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	560089	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	213978	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	404436	20.00	ng	0.00
76) Chrysene-d12	13.87	240	435091	20.00	ng	0.00
87) Perylene-d12	15.29	264	322277	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	207646	20.54	ng	0.00
7) Phenol-d6	6.36	99	263800	20.77	ng	0.00
23) Nitrobenzene-d5	7.30	82	139543	14.91	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	53566	23.58	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	239733	14.14	ng	0.00
79) Terphenyl-d14	12.83	244	250304	11.15	ng	0.00
Target Compounds						
71) Phenanthrene	11.27	178	62910	3.126	ng	96
75) Fluoranthene	12.46	202	173646	8.055	ng	96
78) Pyrene	12.68	202	165375	4.509	ng	98
81) Benzo(a)anthracene	13.86	228	93089	3.087	ng	99
83) Chrysene	13.90	228	90186	3.053	ng	96
88) Benzo(b)fluoranthene	14.89	252	111985m	5.372	ng	
90) Benzo(a)pyrene	15.22	252	62869	3.410	ng	98
92) Benzo(g,h,i)perylene	17.05	276	36279	2.296	ng	94

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

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