

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
 Data File : BF112598.D
 Acq On : 18 Feb 2019 13:57
 Operator : JU/SJ
 Sample : K1358-03 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-021419

Manual Integrations
 APPROVED

Sohil
 2/19/2019 9:38:37 AM

Quant Time: Feb 18 14:39:14 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	120267	20.00	ng	-0.02
21) Naphthalene-d8	8.10	136	477388	20.00	ng	-0.02
39) Acenaphthene-d10	9.86	164	217511	20.00	ng	-0.02
64) Phenanthrene-d10	11.35	188	348638	20.00	ng	-0.02
76) Chrysene-d12	14.00	240	257759	20.00	ng	0.00
87) Perylene-d12	15.48	264	212609	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	331395	58.53	ng	-0.02
7) Phenol-d6	6.47	99	418113	53.14	ng	-0.02
23) Nitrobenzene-d5	7.39	82	268794	32.44	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	112211	44.32	ng	-0.02
45) 2-Fluorobiphenyl	9.17	172	475804	30.94	ng	-0.03
79) Terphenyl-d14	12.93	244	362074	26.46	ng	-0.02
Target Compounds						
71) Phenanthrene	11.37	178	40632	2.242	ng	97
75) Fluoranthene	12.57	202	45692	2.350	ng	98
78) Pyrene	12.81	202	51060	2.861	ng	98
88) Benzo(b)fluoranthene	15.05	252	29651m	2.332	ng	

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

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