

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122818\
 Data File : BF111791.D
 Acq On : 28 Dec 2018 16:23
 Operator : JU/SJ
 Sample : J6193-11 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-122718-A

Manual Integrations
 APPROVED

Sohil
 1/2/2019 2:51:34 PM

Quant Time: Dec 29 01:38:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	161471	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	593646	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	269677	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	420574	20.00	ng	0.01
76) Chrysene-d12	14.07	240	362130	20.00	ng	0.02
87) Perylene-d12	15.56	264	277181	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	196651	20.76	ng	0.02
7) Phenol-d6	6.54	99	247824	20.56	ng	0.02
23) Nitrobenzene-d5	7.46	82	152603	14.96	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	55176	17.32	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	275980	14.92	ng	0.00
79) Terphenyl-d14	13.00	244	219335	11.49	ng	0.00
Target Compounds						
71) Phenanthrene	11.45	178	58008	2.601	ng	96
75) Fluoranthene	12.64	202	89335	3.955	ng	94
78) Pyrene	12.87	202	85934	3.360	ng	95
83) Chrysene	14.09	228	42522	2.093	ng	96
88) Benzo(b)fluoranthene	15.12	252	45806m	2.828	ng	
90) Benzo(a)pyrene	15.49	252	31746	2.157	ng	# 91

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

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