

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111606.D
 Acq On : 19 Dec 2018 22:39
 Operator : JU/SJ
 Sample : J6322-21
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-6-C-2-120718

Manual Integrations
 APPROVED

Sohil
 12/20/2018 12:42:13 PM

Quant Time: Dec 20 04:10:10 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	6.89	152	194890	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	728603	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	366195	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	645082	20.00	ng	0.00
76) Chrysene-d12	14.04	240	518829	20.00	ng	-0.01
87) Perylene-d12	15.52	264	489596	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	1320306	115.48	ng	0.00
7) Phenol-d6	6.52	99	1695048	116.49	ng	0.00
23) Nitrobenzene-d5	7.45	82	1045985	83.56	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	469900	108.60	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1791595	71.31	ng	0.00
79) Terphenyl-d14	12.99	244	1564097	57.17	ng	0.00
Target Compounds						
10) Phenol	6.53	94	100420	6.116	ng	94
50) Dimethylphthalate	9.64	163	142468	5.320	ng	97
71) Phenanthrene	11.43	178	123809	3.619	ng	93
75) Fluoranthene	12.62	202	163414	4.717	ng	95
78) Pyrene	12.85	202	132510	3.617	ng	95
81) Benzo(a)anthracene	14.03	228	70159	2.369	ng	99
83) Chrysene	14.07	228	64365	2.212	ng	96
88) Benzo(b)fluoranthene	15.08	252	75513m	2.639	ng	

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

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