

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111616.D
 Acq On : 20 Dec 2018 3:15
 Operator : JU/SJ
 Sample : J6386-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS006-1824-02

Quant Time: Dec 20 04:46:22 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	197772	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	686983	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	311397	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	524119	20.00	ng	0.00
76) Chrysene-d12	14.05	240	452473	20.00	ng	0.00
87) Perylene-d12	15.53	264	309341	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	1524032	131.35	ng	0.00
7) Phenol-d6	6.53	99	1918351	129.92	ng	0.00
23) Nitrobenzene-d5	7.46	82	1142083	96.77	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	458330	124.57	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1980705	92.71	ng	0.00
79) Terphenyl-d14	13.00	244	1804446	75.63	ng	0.00
Target Compounds						
10) Phenol	6.53	94	138805	8.331	ng	90
50) Dimethylphthalate	9.64	163	221274	9.717	ng	99
71) Phenanthrene	11.44	178	116958	4.208	ng	94
75) Fluoranthene	12.62	202	80950	2.876	ng	95
78) Pyrene	12.85	202	141642	4.433	ng	94
90) Benzo(a)pyrene	15.46	252	35439	2.158	ng	# 69
92) Benzo(g,h,i)perylene	17.45	276	31604	2.250	ng	# 91

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

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