

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111604.D
 Acq On : 19 Dec 2018 21:43
 Operator : JU/SJ
 Sample : J6464-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 290-RT4008-RBR251686

Manual Integrations
 APPROVED

Sohil
 12/20/2018 12:42:10 PM

Quant Time: Dec 20 04:04:23 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	199333	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	754033	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	385843	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	732507	20.00	ng	0.00
76) Chrysene-d12	14.04	240	502537	20.00	ng	-0.01
87) Perylene-d12	15.52	264	493688	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	1432440	122.49	ng	0.00
7) Phenol-d6	6.52	99	1875469	126.02	ng	0.00
23) Nitrobenzene-d5	7.45	82	1148716	88.68	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	561697	123.21	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1957328	73.94	ng	0.00
79) Terphenyl-d14	13.00	244	1804213	68.09	ng	0.00
Target Compounds						
10) Phenol	6.53	94	98709	5.878	ng	90
50) Dimethylphthalate	9.64	163	161044	5.708	ng	# 97
71) Phenanthrene	11.43	178	107794	2.775	ng	94
75) Fluoranthene	12.62	202	179711	4.568	ng	96
78) Pyrene	12.85	202	158456	4.465	ng	95
81) Benzo(a)anthracene	14.03	228	75355	2.627	ng	93
83) Chrysene	14.07	228	78342	2.779	ng	97
88) Benzo(b)fluoranthene	15.08	252	121226m	4.201	ng	
90) Benzo(a)pyrene	15.44	252	70532m	2.691	ng	

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

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