

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031219\
 Data File : BF113018.D
 Acq On : 12 Mar 2019 22:27
 Operator : JU/SJ
 Sample : K1847-17
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-7B-C-4-031119

Manual Integrations
 APPROVED

Sohil
 3/13/2019 11:40:46 AM

Quant Time: Mar 13 04:44:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	179258	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	681206	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	317305	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	574980	20.00	ng	0.00
76) Chrysene-d12	13.86	240	481879	20.00	ng	0.00
87) Perylene-d12	15.27	264	408895	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1166296	101.21	ng	0.00
7) Phenol-d6	6.37	99	1508793	104.23	ng	0.00
23) Nitrobenzene-d5	7.29	82	938751	82.46	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	435104	129.17	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1648029	86.44	ng	0.00
79) Terphenyl-d14	12.82	244	1499040	60.30	ng	0.00
Target Compounds						
10) Phenol	6.38	94	42707	2.528	ng	94
50) Dimethylphthalate	9.48	163	151236	6.464	ng	99
71) Phenanthrene	11.26	178	213889	7.477	ng	96
75) Fluoranthene	12.45	202	488758	15.947	ng	97
78) Pyrene	12.67	202	433449	10.670	ng	99
81) Benzo(a)anthracene	13.86	228	205412	6.151	ng	98
83) Chrysene	13.89	228	208601	6.377	ng	97
86) Indeno(1,2,3-cd)pyrene	16.63	276	77716	2.787	ng	# 93
88) Benzo(b)fluoranthene	14.87	252	230299m	8.707	ng	
89) Benzo(k)fluoranthene	14.89	252	93373m	3.898	ng	
90) Benzo(a)pyrene	15.21	252	160070	6.842	ng	98
92) Benzo(a,h,i)perylene	17.03	276	76267	3.805	ng	99

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

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