

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012419\
 Data File : BF112219.D
 Acq On : 24 Jan 2019 20:01
 Operator : JU/SJ
 Sample : K1223-07 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-S

Manual Integrations
 APPROVED

Sohil
 1/25/2019 10:05:21 AM

Quant Time: Jan 25 00:42:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 19 01:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	211480	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	784582	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	281772	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	567078	20.00	ng	0.00
76) Chrysene-d12	14.03	240	425673	20.00	ng	0.00
87) Perylene-d12	15.51	264	310303	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	556504	47.76	ng	0.00
7) Phenol-d6	6.50	99	688456	47.23	ng	0.00
23) Nitrobenzene-d5	7.42	82	397056	34.96	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	168756	55.52	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	593365	30.87	ng	0.00
79) Terphenyl-d14	12.96	244	602144	30.09	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.60	163	67288	3.275	ng	99
71) Phenanthrene	11.40	178	121771	4.267	ng	98
75) Fluoranthene	12.60	202	187036	6.207	ng	97
78) Pyrene	12.83	202	161373	5.746	ng	99
80) Butylbenzylphthalate	13.43	149	25849	2.098	ng	# 87
81) Benzo(a)anthracene	14.02	228	72424	2.967	ng	96
83) Chrysene	14.06	228	71134	3.080	ng	97
88) Benzo(b)fluoranthene	15.07	252	72337m	4.107	ng	
90) Benzo(a)pyrene	15.44	252	44134	2.732	ng	# 69

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

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