

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
 Data File : BF111989.D
 Acq On : 10 Jan 2019 16:06
 Operator : JU/SJ
 Sample : K1015-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-010918-A

Manual Integrations
 APPROVED

Sohil
 1/11/2019 4:42:52 PM

Quant Time: Jan 11 01:17:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	117251	20.00	ng	-0.02
21) Naphthalene-d8	8.16	136	442231	20.00	ng	-0.02
39) Acenaphthene-d10	9.91	164	184353	20.00	ng	-0.02
64) Phenanthrene-d10	11.40	188	345671	20.00	ng	-0.02
76) Chrysene-d12	14.04	240	277079	20.00	ng	-0.01
87) Perylene-d12	15.53	264	237949	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	725823	105.91	ng	-0.01
7) Phenol-d6	6.52	99	847167	95.91	ng	-0.02
23) Nitrobenzene-d5	7.43	82	514091	62.13	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	224965	93.82	ng	-0.02
45) 2-Fluorobiphenyl	9.23	172	818663	64.75	ng	-0.02
79) Terphenyl-d14	12.99	244	641925	43.98	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	9.62	163	161993	11.697	ng	99
71) Phenanthrene	11.42	178	357564	19.347	ng	99
72) Anthracene	11.47	178	99712	5.315	ng	99
75) Fluoranthene	12.62	202	589165	31.053	ng	99
78) Pyrene	12.84	202	452166	23.717	ng	98
81) Benzo(a)anthracene	14.03	228	247737	15.536	ng	99
83) Chrysene	14.07	228	212760	13.993	ng	96
86) Indeno(1,2,3-cd)pyrene	17.01	276	71126	6.417	ng	96
88) Benzo(b)fluoranthene	15.09	252	233914m	16.434	ng	
89) Benzo(k)fluoranthene	15.12	252	68957m	5.154	ng	
90) Benzo(a)pyrene	15.46	252	147932	11.820	ng	100
92) Benzo(a,h,i)perylene	17.46	276	61453	6.242	ng	97

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

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