

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
 Data File : BF112003.D
 Acq On : 10 Jan 2019 22:27
 Operator : JU/SJ
 Sample : K1015-03 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 1/11/2019 4:42:59 PM

Quant Time: Jan 11 02:48:36 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	87532	20.00	ng	-0.02
21) Naphthalene-d8	8.16	136	299732	20.00	ng	-0.02
39) Acenaphthene-d10	9.92	164	138199	20.00	ng	-0.02
64) Phenanthrene-d10	11.40	188	275811	20.00	ng	-0.02
76) Chrysene-d12	14.04	240	248761	20.00	ng	-0.02
87) Perylene-d12	15.52	264	186054	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	264157	51.63	ng	-0.02
7) Phenol-d6	6.51	99	323423	49.04	ng	-0.02
23) Nitrobenzene-d5	7.43	82	185178	33.02	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	84465	46.99	ng	-0.02
45) 2-Fluorobiphenyl	9.23	172	301448	31.81	ng	-0.02
79) Terphenyl-d14	12.98	244	382089	29.16	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.62	163	67208	6.474	ng	Qvalue 98
52) Acenaphthene	9.95	154	19896	2.301	ng	99
58) Fluorene	10.46	166	27865	3.045	ng	96
71) Phenanthrene	11.42	178	276340	18.740	ng	99
72) Anthracene	11.47	178	70127	4.685	ng	96
75) Fluoranthene	12.62	202	390768	25.813	ng	99
78) Pyrene	12.84	202	353958	20.680	ng	99
81) Benzo(a)anthracene	14.03	228	148091	10.344	ng	98
83) Chrysene	14.07	228	121509	8.901	ng	97
86) Indeno(1,2,3-cd)pyrene	17.01	276	58859	5.914	ng	95
88) Benzo(b)fluoranthene	15.09	252	125348m	11.263	ng	
89) Benzo(k)fluoranthene	15.12	252	46611m	4.455	ng	
90) Benzo(a)pyrene	15.46	252	98422	10.057	ng	# 96
92) Benzo(a,h,i)perylene	17.46	276	56162	7.295	ng	96

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

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