

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010819\
 Data File : BF111930.D
 Acq On : 9 Jan 2019 1:45
 Operator : JU/SJ
 Sample : K1044-07 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPSB-09(10-15)

Quant Time: Jan 09 03:12:29 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	6.89	152	115813	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	372048	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	188887	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	285000	20.00	ng	0.00
76) Chrysene-d12	14.07	240	185298	20.00	ng	0.02
87) Perylene-d12	15.58	264	158188	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	87464	12.92	ng	0.00
7) Phenol-d6	6.53	99	99881	11.45	ng	0.00
23) Nitrobenzene-d5	7.44	82	51952	7.46	ng	-0.02
42) 2,4,6-Tribromophenol	10.72	330	22831	9.29	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	93772	7.24	ng	-0.01
79) Terphenyl-d14	13.00	244	82369	8.44	ng	0.00
Target Compounds						
52) Acenaphthene	9.96	154	32626	2.760	ng	98
58) Fluorene	10.47	166	25882	2.069	ng	# 89
71) Phenanthrene	11.44	178	35878	2.355	ng	# 80
75) Fluoranthene	12.64	202	107144	6.849	ng	93
78) Pyrene	12.87	202	99343	7.792	ng	# 93
81) Benzo(a)anthracene	14.06	228	34237	3.210	ng	# 64
83) Chrysene	14.10	228	39268	3.862	ng	# 63

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

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