

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
 Data File : BF106985.D
 Acq On : 29 Jun 2018 20:40
 Operator : JU/SJ
 Sample : J3692-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-02

Quant Time: Jun 30 00:22:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	73603	20.00	ng	-0.01
21) Naphthalene-d8	8.24	136	266852	20.00	ng	-0.01
38) Acenaphthene-d10	9.99	164	119964	20.00	ng	-0.02
63) Phenanthrene-d10	11.48	188	200042	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	195415	20.00	ng	-0.03
86) Perylene-d12	15.67	264	145165	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	255947	58.88	ng	-0.01
7) Phenol-d6	6.59	99	196338	36.17	ng	-0.02
23) Nitrobenzene-d5	7.52	82	418532	87.16	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	183637	132.07	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	738421	108.22	ng	-0.02
78) Terphenyl-d14	13.07	244	735160	91.13	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.61	94	17360	2.802	ng	97
31) Naphthalene	8.26	128	463494	37.150	ng	99
37) 2-Methylnaphthalene	8.95	142	34768	4.204	ng	93
51) Acenaphthene	10.02	154	46825	6.259	ng	95
54) Dibenzofuran	10.20	168	41234	4.025	ng	95
57) Fluorene	10.54	166	37307	4.589	ng	99
70) Phenanthrene	11.51	178	68190	7.067	ng	97
72) Carbazole	11.72	167	29044	3.150	ng	98

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

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