

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090717\
 Data File : BF098338.D
 Acq On : 8 Sep 2017 3:15
 Operator : SJ/JU
 Sample : I5131-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 090517-TIDO-F-D-B

Quant Time: Sep 08 04:38:04 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	103423	20.00	ng	-0.02
21) Naphthalene-d8	7.97	136	397608	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	165844	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	300218	20.00	ng	-0.01
75) Chrysene-d12	13.84	240	239723	20.00	ng	-0.02
86) Perylene-d12	15.25	264	166554	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	328729	48.35	ng	-0.01
7) Phenol-d6	6.33	99	285914	30.95	ng	-0.02
23) Nitrobenzene-d5	7.26	82	713719	97.51	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	137867	95.81	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	1064714	94.32	ng	-0.01
78) Terphenyl-d14	12.80	244	832359	72.41	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.00	128	458379	22.206	ng	99
37) 2-Methylnaphthalene	8.69	142	75790	5.989	ng	98
49) Dimethylphthalate	9.44	163	25445	2.099	ng	98
51) Acenaphthene	9.76	154	106256	9.601	ng	97
54) Dibenzofuran	9.93	168	112251	7.568	ng	99
57) Fluorene	10.27	166	95713	8.346	ng	98
70) Phenanthrene	11.23	178	276888	17.308	ng	98
72) Carbazole	11.44	167	31016	2.014	ng	# 96
74) Fluoranthene	12.42	202	100901	6.043	ng	97
77) Pyrene	12.64	202	64428	3.286	ng	97

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

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