

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110916\
 Data File : BF091135.D
 Acq On : 9 Nov 2016 20:18
 Operator : UM/SJ
 Sample : H5602-01DL 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 USTD

Quant Time: Nov 09 23:16:04 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 12:59:29 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	200373	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	768280	20.00	ng	-0.01
38) Acenaphthene-d10	9.69	164	478499	20.00	ng	-0.01
63) Phenanthrene-d10	11.17	188	719683	20.00	ng	0.00
75) Chrysene-d12	13.80	240	471345	20.00	ng	-0.01
86) Perylene-d12	15.17	264	366016	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	2369	0.20	ng	0.15
7) Phenol-d6	6.34	99	18255	1.11	ng	0.02
23) Nitrobenzene-d5	7.22	82	609120	43.25	ng	-0.02
41) 2,4,6-Tribromophenol	10.49	330	246508	56.98	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1143986	41.16	ng	-0.01
78) Terphenyl-d14	12.75	244	790434	41.85	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	7.96	128	258981	7.05	ng	# 95
37) 2-Methylnaphthalene	8.66	142	1108014	46.91	ng	100
45) 1,1'-Biphenyl	9.12	154	116276	3.33	ng	89
51) Acenaphthene	9.72	154	56994	2.20	ng	97
54) Dibenzofuran	9.89	168	79185	2.27	ng	# 87
57) Fluorene	10.24	166	105855	3.71	ng	# 96
70) Phenanthrene	11.20	178	134902	3.53	ng	98

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

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