

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
 Data File : BF091182.D
 Acq On : 15 Nov 2016 16:49
 Operator : UM/SJ
 Sample : H5636-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPS02304

Quant Time: Nov 16 06:34:35 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	6.64	152	165626	20.00	ng	-0.02
21) Naphthalene-d8	7.92	136	705428	20.00	ng	-0.03
38) Acenaphthene-d10	9.68	164	372297	20.00	ng	-0.02
63) Phenanthrene-d10	11.15	188	662920	20.00	ng	-0.02
75) Chrysene-d12	13.78	240	591633	20.00	ng	-0.03
86) Perylene-d12	15.14	264	437185	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	661341	65.95	ng	-0.02
7) Phenol-d6	6.28	99	558898	41.06	ng	-0.03
23) Nitrobenzene-d5	7.21	82	1212245	93.74	ng	-0.03
41) 2,4,6-Tribromophenol	10.46	330	536276	159.33	ng	-0.02
44) 2-Fluorobiphenyl	9.00	172	1834910	84.85	ng	-0.03
78) Terphenyl-d14	12.74	244	2142408	90.36	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	7.94	128	546633	16.20	ng	98
37) 2-Methylnaphthalene	8.64	142	133320	6.15	ng	98
49) Dimethylphthalate	9.40	163	105102	4.31	ng	99
51) Acenaphthene	9.70	154	91851	4.55	ng	99
54) Dibenzofuran	9.88	168	140479	5.17	ng	# 85
57) Fluorene	10.21	166	178302	8.02	ng	# 95
70) Phenanthrene	11.17	178	250346	7.10	ng	97

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

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