

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072417\
 Data File : BF097011.D
 Acq On : 25 Jul 2017 3:17
 Operator : SJ/JU
 Sample : I4360-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CEMETERY-11-(0-2)

Quant Time: Jul 25 04:01:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 01:37:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.52	152	135961	20.00	ng	-0.02
21) Naphthalene-d8	7.80	136	542648	20.00	ng	-0.03
38) Acenaphthene-d10	9.54	164	219006	20.00	ng	-0.04
63) Phenanthrene-d10	11.02	188	321784	20.00	ng	-0.04
75) Chrysene-d12	13.64	240	246252	20.00	ng	-0.04
86) Perylene-d12	15.00	264	220664	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	760149	84.06	ng	-0.03
7) Phenol-d6	6.18	99	981456	90.45	ng	-0.04
23) Nitrobenzene-d5	7.08	82	589892	67.33	ng	-0.04
41) 2,4,6-Tribromophenol	10.33	330	154710	82.76	ng	-0.04
44) 2-Fluorobiphenyl	8.87	172	904370	65.24	ng	-0.04
78) Terphenyl-d14	12.60	244	542376	38.20	ng	-0.04
Target Compounds						
10) Phenol	6.19	94	32830	2.676	ng	100
49) Dimethylphthalate	9.27	163	118869	7.208	ng	99
85) Indeno(1,2,3-cd)pyrene	16.23	276	8252	3.897	ng	# 84

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

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