

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081018\
 Data File : BG036280.D
 Acq On : 10 Aug 2018 23:20
 Operator : JU/SJ
 Sample : J4272-22
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JC-03-080918-K

Quant Time: Aug 11 04:37:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 31 12:30:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	27153	20.00	ng	-0.02
21) Naphthalene-d8	10.81	136	94560	20.00	ng	-0.02
38) Acenaphthene-d10	14.63	164	65339	20.00	ng	-0.02
63) Phenanthrene-d10	17.38	188	179480	20.00	ng	-0.01
75) Chrysene-d12	21.64	240	193498	20.00	ng	-0.02
86) Perylene-d12	24.84	264	167092	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	229194	140.14	ng	0.00
7) Phenol-d6	7.18	99	288825	118.36	ng	-0.01
23) Nitrobenzene-d5	9.17	82	256993	113.53	ng	-0.02
41) 2,4,6-Tribromophenol	16.12	330	142165	165.08	ng	-0.02
44) 2-Fluorobiphenyl	13.26	172	592236	121.43	ng	-0.02
78) Terphenyl-d14	19.98	244	1095654	124.82	ng	-0.02
Target Compounds						
19) n-Nitroso-di-n-propylamine	9.17	70	35916	17.479	ng	# 64
50) 2,6-Dinitrotoluene	14.63	165	9049	7.762	ng	# 21
56) 2,4-Dinitrotoluene	14.63	165	9049	5.411	ng	# 19
76) Benzidine	19.98	184	19154	3.765	ng	# 87

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

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