

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080918\
 Data File : BG036240.D
 Acq On : 9 Aug 2018 19:35
 Operator : JU/SJ
 Sample : J4379-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-103S

Quant Time: Aug 10 05:13:31 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	30450	20.00	ng	-0.01
21) Naphthalene-d8	10.81	136	113335	20.00	ng	-0.02
38) Acenaphthene-d10	14.63	164	76525	20.00	ng	-0.02
63) Phenanthrene-d10	17.37	188	191091	20.00	ng	-0.02
75) Chrysene-d12	21.64	240	206568	20.00	ng	-0.02
86) Perylene-d12	24.82	264	202590	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	111389	60.73	ng	0.00
7) Phenol-d6	7.18	99	99419	36.33	ng	-0.01
23) Nitrobenzene-d5	9.17	82	242460	89.37	ng	-0.01
41) 2,4,6-Tribromophenol	16.12	330	158111	156.76	ng	-0.02
44) 2-Fluorobiphenyl	13.26	172	525193	91.95	ng	-0.02
78) Terphenyl-d14	19.99	244	878443	93.74	ng	-0.01
Target Compounds						
31) Naphthalene	10.87	128	24743	4.191	ng	# 94
49) Dimethylphthalate	14.08	163	14059	2.097	ng	# 96
51) Acenaphthene	14.70	154	26120	5.424	ng	97
70) Phenanthrene	17.42	178	19733	2.070	ng	# 92
74) Fluoranthene	19.43	202	35994	2.802	ng	97

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

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