

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080218\
 Data File : BG036064.D
 Acq On : 2 Aug 2018 17:28
 Operator : JU/SJ
 Sample : J4262-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GW-01

Quant Time: Aug 02 18:55:08 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	8.02	152	29363	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	115031	20.00	ng	0.00
38) Acenaphthene-d10	14.64	164	83270	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	234000	20.00	ng	0.00
75) Chrysene-d12	21.65	240	262971	20.00	ng	0.00
86) Perylene-d12	24.84	264	258920	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	95867	54.20	ng	0.00
7) Phenol-d6	7.19	99	83593	31.68	ng	0.00
23) Nitrobenzene-d5	9.18	82	218024	79.18	ng	0.00
41) 2,4,6-Tribromophenol	16.13	330	161967	147.57	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	534598	86.01	ng	0.00
78) Terphenyl-d14	19.99	244	1076883	90.27	ng	0.00
Target Compounds						
31) Naphthalene	10.88	128	14108	2.354	ng	# 93
49) Dimethylphthalate	14.10	163	30637	4.199	ng	# 95
51) Acenaphthene	14.71	154	11702	2.233	ng	# 83
70) Phenanthrene	17.43	178	47473	4.066	ng	97

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

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