

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
 Data File : BF107927.D
 Acq On : 2 Aug 2018 23:23
 Operator : JU/SJ
 Sample : J4262-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-03

Quant Time: Aug 03 04:54:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	137522	20.00	ng	-0.01
21) Naphthalene-d8	8.24	136	556990	20.00	ng	-0.01
38) Acenaphthene-d10	9.99	164	221806	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	419504	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	366190	20.00	ng	-0.01
86) Perylene-d12	15.67	264	309627	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	415752	51.38	ng	0.00
7) Phenol-d6	6.61	99	324205	30.40	ng	0.00
23) Nitrobenzene-d5	7.52	82	872294	90.22	ng	-0.01
41) 2,4,6-Tribromophenol	10.79	330	340815	113.08	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	1572610	98.80	ng	-0.01
78) Terphenyl-d14	13.07	244	1441129	85.34	ng	-0.01
Target Compounds						
31) Naphthalene	8.26	128	488218	18.856	ng	99
32) Benzoic acid	8.05	122	11800	3.994	ng	# 56
37) 2-Methylnaphthalene	8.95	142	62756	3.859	ng	98
49) Dimethylphthalate	9.70	163	57855	3.473	ng	99
51) Acenaphthene	10.02	154	35509	2.590	ng	97
57) Fluorene	10.55	166	38335	2.328	ng	100
70) Phenanthrene	11.52	178	43698	2.194	ng	97

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

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