

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080718\
 Data File : BG036171.D
 Acq On : 7 Aug 2018 21:01
 Operator : JU/SJ
 Sample : J4366-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SV30518L

Quant Time: Aug 08 02:03:58 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	30176	20.00	ng	-0.01
21) Naphthalene-d8	10.81	136	106035	20.00	ng	-0.02
38) Acenaphthene-d10	14.64	164	75375	20.00	ng	-0.01
63) Phenanthrene-d10	17.37	188	216584	20.00	ng	-0.02
75) Chrysene-d12	21.64	240	252120	20.00	ng	-0.02
86) Perylene-d12	24.82	264	241821	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	132374	72.83	ng	0.00
7) Phenol-d6	7.18	99	121016	44.63	ng	-0.02
23) Nitrobenzene-d5	9.17	82	253530	99.88	ng	-0.01
41) 2,4,6-Tribromophenol	16.12	330	182608	183.80	ng	-0.02
44) 2-Fluorobiphenyl	13.26	172	587390	104.40	ng	-0.02
78) Terphenyl-d14	19.99	244	1234183	107.91	ng	-0.01
Target Compounds						
49) Dimethylphthalate	14.09	163	17130	2.594	ng	98
51) Acenaphthene	14.70	154	45601	9.614	ng	95
54) Dibenzofuran	15.04	168	26631	3.530	ng	92
57) Fluorene	15.68	166	38331	6.063	ng	99
70) Phenanthrene	17.42	178	49921	4.619	ng	98

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

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