

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
 Data File : BG036222.D
 Acq On : 9 Aug 2018 7:51
 Operator : JU/SJ
 Sample : J4328-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-22

Quant Time: Aug 09 09:03:31 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	29580	20.00	ng	-0.01
21) Naphthalene-d8	10.81	136	103613	20.00	ng	-0.02
38) Acenaphthene-d10	14.63	164	77906	20.00	ng	-0.02
63) Phenanthrene-d10	17.37	188	205697	20.00	ng	-0.02
75) Chrysene-d12	21.64	240	228403	20.00	ng	-0.02
86) Perylene-d12	24.83	264	215273	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	94781	53.20	ng	0.00
7) Phenol-d6	7.18	99	82260	30.95	ng	-0.01
23) Nitrobenzene-d5	9.17	82	229922	92.70	ng	-0.02
41) 2,4,6-Tribromophenol	16.12	330	160061	155.88	ng	-0.02
44) 2-Fluorobiphenyl	13.26	172	512025	88.05	ng	-0.02
78) Terphenyl-d14	19.99	244	1006516	97.14	ng	-0.01
Target Compounds						
51) Acenaphthene	14.70	154	49990	10.197	ng	91
54) Dibenzofuran	15.03	168	49497	6.348	ng	# 89
57) Fluorene	15.68	166	16232	2.484	ng	# 91

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

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