

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG010218\
 Data File : BG031875.D
 Acq On : 2 Jan 2018 17:25
 Operator : SJ/JU
 Sample : I6680-24
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CL-01-122917-B

Manual Integrations
 APPROVED

Sohil
 1/3/2018 10:25:16 AM

Quant Time: Jan 03 00:19:26 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 29 15:51:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.79	152	46984	20.00	ng	0.00
21) Naphthalene-d8	11.67	136	192122	20.00	ng	0.00
38) Acenaphthene-d10	15.41	164	140292	20.00	ng	-0.01
63) Phenanthrene-d10	18.15	188	396826	20.00	ng	-0.01
75) Chrysene-d12	22.49	240	470729	20.00	ng	-0.02
86) Perylene-d12	26.22	264	487331	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	354236	137.32	ng	0.00
7) Phenol-d6	7.89	99	413439	123.45	ng	0.00
23) Nitrobenzene-d5	9.99	82	278252	109.36	ng	0.00
41) 2,4,6-Tribromophenol	16.89	330	387368	180.96	ng	0.00
44) 2-Fluorobiphenyl	14.05	172	999433	89.42	ng	0.00
78) Terphenyl-d14	20.68	244	2089182	101.40	ng	-0.01
Target Compounds						
31) Naphthalene	11.73	128	688032	70.961	ng	99
37) 2-Methylnaphthalene	13.28	142	37319m	4.748	ng	
48) Acenaphthylene	15.14	152	54767	3.746	ng	99
54) Dibenzofuran	15.81	168	32565	2.218	ng	# 96
57) Fluorene	16.45	166	30923	2.593	ng	# 93
70) Phenanthrene	18.19	178	56496	2.516	ng	98

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

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