

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012318\
 Data File : BG032231.D
 Acq On : 23 Jan 2018 14:44
 Operator : SJ/JU
 Sample : J1239-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-7317-01

Quant Time: Jan 24 04:23:27 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 18 10:13:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.73	152	47501	20.00	ng	-0.01
21) Naphthalene-d8	11.61	136	198262	20.00	ng	-0.01
38) Acenaphthene-d10	15.36	164	138918	20.00	ng	-0.02
63) Phenanthrene-d10	18.09	188	341010	20.00	ng	-0.02
75) Chrysene-d12	22.43	240	448509	20.00	ng	-0.02
86) Perylene-d12	26.12	264	469678	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	6.16	112	287079	110.53	ng	0.00
7) Phenol-d6	7.84	99	336100	102.75	ng	0.00
23) Nitrobenzene-d5	9.93	82	251323	85.76	ng	-0.01
41) 2,4,6-Tribromophenol	16.84	330	275628	119.90	ng	-0.01
44) 2-Fluorobiphenyl	13.99	172	860100	76.77	ng	-0.01
78) Terphenyl-d14	20.64	244	1690885	83.24	ng	-0.01
Target Compounds						
10) Phenol	7.87	94	11127	3.453	ng	Qvalue 92
31) Naphthalene	11.67	128	2763357	281.360	ng	# 94
37) 2-Methylnaphthalene	13.22	142	347560	44.574	ng	96
45) 1,1'-Biphenyl	14.20	154	58764	4.934	ng	93
51) Acenaphthene	15.43	154	154982	16.613	ng	98
57) Fluorene	16.40	166	88067	7.716	ng	95
70) Phenanthrene	18.14	178	204342	10.763	ng	99
72) Carbazole	18.49	167	60845	3.704	ng	97

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

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