

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032918\
 Data File : BG033446.D
 Acq On : 30 Mar 2018 9:23
 Operator : SJ/JU
 Sample : J2102-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 7005

Quant Time: Mar 30 10:59:53 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 28 18:34:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.87	152	32944	20.00	ng	0.00
21) Naphthalene-d8	11.75	136	153748	20.00	ng	0.00
38) Acenaphthene-d10	15.48	164	113817	20.00	ng	0.00
63) Phenanthrene-d10	18.21	188	299423	20.00	ng	0.00
75) Chrysene-d12	22.56	240	321253	20.00	ng	0.00
86) Perylene-d12	26.33	264	343643	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.30	112	282251	160.65	ng	0.00
7) Phenol-d6	7.98	99	433114	143.48	ng	0.00
23) Nitrobenzene-d5	10.07	82	347424	103.82	ng	0.00
41) 2,4,6-Tribromophenol	16.96	330	274803	161.43	ng	0.00
44) 2-Fluorobiphenyl	14.11	172	829244	105.18	ng	0.00
78) Terphenyl-d14	20.73	244	1287216	85.69	ng	0.00
Target Compounds						
10) Phenol	8.02	94	11417	3.831	ng	# 61
15) Benzyl Alcohol	9.10	79	75073	31.587	ng	# 85
20) 3+4-Methylphenols	9.64	107	41057	14.203	ng	88
31) Naphthalene	11.80	128	22897	2.874	ng	97
32) Benzoic acid	11.11	122	240833	131.508	ng	94

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

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