

Data Path : Z:\HPCHEM1\BNA G\DATA\BG040918\
 Data File : BG033682.D
 Acq On : 9 Apr 2018 16:01
 Operator : SJ/JU
 Sample : J2236-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-103

Quant Time: Apr 10 07:00:17 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 05 17:36:07 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.86	152	42647	20.00	ng	0.00
21) Naphthalene-d8	11.74	136	177995	20.00	ng	0.00
38) Acenaphthene-d10	15.47	164	136978	20.00	ng	0.00
63) Phenanthrene-d10	18.21	188	352553	20.00	ng	0.00
75) Chrysene-d12	22.55	240	343016	20.00	ng	-0.01
86) Perylene-d12	26.31	264	368939	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.29	112	140982	61.99	ng	0.00
7) Phenol-d6	7.97	99	129885	33.24	ng	0.00
23) Nitrobenzene-d5	10.06	82	329095	84.95	ng	0.00
41) 2,4,6-Tribromophenol	16.95	330	242527	118.38	ng	0.00
44) 2-Fluorobiphenyl	14.10	172	787346	82.98	ng	0.00
78) Terphenyl-d14	20.73	244	1369368	85.38	ng	0.00
Target Compounds						
3) Pyridine	4.39	79	23780	7.448	ng	94
10) Phenol	8.01	94	8631	2.237	ng	78
20) 3+4-Methylphenols	9.63	107	10150	2.712	ng	83

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

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