

Data Path : Z:\HPCHEM1\BNA M\DATA\BM052617\
 Data File : BM010251.D
 Acq On : 26 May 2017 11:35
 Operator : SJ/MA
 Sample : I3340-13
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DRUM-1-DRUM-2-COMP

Quant Time: May 27 00:11:33 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051917.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 20 04:25:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.94	152	117702	20.00	ng	-0.04
21) Naphthalene-d8	10.75	136	413998	20.00	ng	-0.04
38) Acenaphthene-d10	14.57	164	290337	20.00	ng	-0.04
63) Phenanthrene-d10	17.32	188	758756	20.00	ng	-0.03
75) Chrysene-d12	21.48	240	1092698	20.00	ng	-0.03
86) Perylene-d12	23.83	264	1150359	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	446609	68.52	ng	-0.03
7) Phenol-d6	7.13	99	372115	44.37	ng	-0.04
23) Nitrobenzene-d5	9.13	82	836786	109.05	ng	-0.04
41) 2,4,6-Tribromophenol	16.07	330	631503	143.21	ng	-0.03
44) 2-Fluorobiphenyl	13.19	172	2190761	97.00	ng	-0.04
78) Terphenyl-d14	19.92	244	4297781	89.43	ng	-0.03
Target Compounds						Qvalue
3) Pyridine	3.85	79	149432	18.64	ng	89
10) Phenol	7.16	94	26367	2.98	ng	91
17) 2-Methylphenol	8.40	107	68658	10.92	ng	# 81
20) 3+4-Methylphenols	8.74	107	18894	2.22	ng	# 79
22) Acetophenone	8.77	105	232432	21.90	ng	# 96
49) Dimethylphthalate	14.03	163	111758	4.32	ng	# 98

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

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