

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041416\
 Data File : BG021670.D
 Acq On : 15 Apr 2016 2:30
 Operator : UM/SJ
 Sample : H2517-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S8-0573

Quant Time: Apr 15 14:26:46 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG041316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 15:48:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.22	152	36414	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	172607	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	108935	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	221630	20.00	ng	0.00
75) Chrysene-d12	21.84	240	249157	20.00	ng	0.00
86) Perylene-d12	25.18	264	244043	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.77	112	159840	73.10	ng	0.01
7) Phenol-d6	7.37	99	149058	45.63	ng	0.01
23) Nitrobenzene-d5	9.39	82	324299	96.57	ng	0.00
41) 2,4,6-Tribromophenol	16.31	330	178761	152.26	ng	0.00
44) 2-Fluorobiphenyl	13.46	172	711952	95.76	ng	0.00
78) Terphenyl-d14	20.15	244	656961	65.79	ng	0.00
Target Compounds						Qvalue
15) Benzyl Alcohol	8.45	79	7995	3.20	ng	# 83
20) 3+4-Methylphenols	8.98	107	13954	4.20	ng	93
22) Acetophenone	9.04	105	169774	35.31	ng	# 96
27) 2,4-Dimethylphenol	10.19	122	35456	11.37	ng	96
31) Naphthalene	11.08	128	651513	70.53	ng	# 93
32) Benzoic acid	10.39	122	35876	24.97	ng	# 83
37) 2-Methylnaphthalene	12.67	142	171716	27.08	ng	97

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

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