

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110918\
 Data File : BG037921.D
 Acq On : 9 Nov 2018 15:24
 Operator : JU/SJ
 Sample : J5860-22
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EP2-SB-110(17-19)

Quant Time: Nov 12 13:15:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 18 14:06:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	52225	20.00	ng	0.00
21) Naphthalene-d8	10.94	136	185559	20.00	ng	0.00
39) Acenaphthene-d10	14.73	164	131949	20.00	ng	-0.01
64) Phenanthrene-d10	17.47	188	281068	20.00	ng	-0.01
76) Chrysene-d12	21.75	240	250920	20.00	ng	-0.02
87) Perylene-d12	25.09	264	238559	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	371760	119.01	ng	0.00
7) Phenol-d6	7.25	99	518878	109.85	ng	0.00
23) Nitrobenzene-d5	9.30	82	349795	105.60	ng	0.02
42) 2,4,6-Tribromophenol	16.21	330	195065	135.54	ng	-0.01
45) 2-Fluorobiphenyl	13.35	172	768491	87.82	ng	-0.01
79) Terphenyl-d14	20.07	244	1073850	91.45	ng	-0.01
Target Compounds						
10) Phenol	7.28	94	44122	8.853	ng	# 62
31) Naphthalene	11.01	128	4897946	537.398	ng	# 72
37) 2-Methylnaphthalene	12.60	142	3995346	601.358	ng	# 80
38) 1-Methylnaphthalene	12.81	142	2086592	323.395	ng	# 98
46) 1,1'-Biphenyl	13.56	154	162070	14.831	ng	99
50) Dimethylphthalate	14.17	163	139164	13.633	ng	99
58) Fluorene	15.77	166	25558	2.690	ng	# 95
71) Phenanthrene	17.52	178	39737	2.782	ng	97

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

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