

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110118\
 Data File : BG037727.D
 Acq On : 1 Nov 2018 20:29
 Operator : JU/SJ
 Sample : J5755-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-2A

Quant Time: Nov 02 07:41:10 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	8.10	152	52095	20.00	ng	0.00
21) Naphthalene-d8	10.92	136	239183	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	146301	20.00	ng	0.00
64) Phenanthrene-d10	17.49	188	277893	20.00	ng	0.00
76) Chrysene-d12	21.77	240	233756	20.00	ng	0.00
87) Perylene-d12	25.11	264	235792	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	178541	57.30	ng	0.00
7) Phenol-d6	7.25	99	169770	36.03	ng	0.00
23) Nitrobenzene-d5	9.28	82	367161	85.99	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	184584	115.68	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	833554	85.92	ng	0.00
79) Terphenyl-d14	20.08	244	965882	88.29	ng	0.00
Target Compounds						
10) Phenol	7.27	94	15113	3.040	ng	91
50) Dimethylphthalate	14.18	163	31806	2.810	ng	# 98
52) Acenaphthene	14.80	154	240165	28.281	ng	99
55) Dibenzofuran	15.14	168	51001	3.625	ng	99
58) Fluorene	15.78	166	21840	2.073	ng	# 93

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

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