

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101220\
 Data File : BG046862.D
 Acq On : 13 Oct 2020 4:28
 Operator : CG/JU
 Sample : L4339-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 QUAD-5

Quant Time: Oct 13 05:56:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 07 11:38:17 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.106	152	57084	20.00	ng	# 0.00
21) Naphthalene-d8	10.915	136	216297	20.00	ng	0.00
39) Acenaphthene-d10	14.728	164	163509	20.00	ng	0.00
64) Phenanthrene-d10	17.466	188	431317	20.00	ng	#-0.01
76) Chrysene-d12	21.743	240	450346	20.00	ng	-0.01
86) Perylene-d12	25.010	264	469040	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.662	112	179790	50.46	ng	0.00
7) Phenol-d6	7.260	99	343428	68.33	ng	0.00
23) Nitrobenzene-d5	9.264	82	289103	71.11	ng	0.00
42) 2,4,6-Tribromophenol	16.208	330	109225	50.66	ng	0.00
45) 2-Fluorobiphenyl	13.353	172	700854	60.68	ng	0.00
79) Terphenyl-d14	20.074	244	1324616	58.95	ng	0.00
Target Compounds						
10) Phenol	7.289	94	10306	2.06	ng	Qvalue 74

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

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