

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101220\
 Data File : BG046869.D
 Acq On : 13 Oct 2020 9:14
 Operator : CG/JU
 Sample : L4339-08
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 QUAD-8

Quant Time: Oct 13 11:06:29 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	54525	20.00	ng	# 0.00
21) Naphthalene-d8	10.914	136	207293	20.00	ng	0.00
39) Acenaphthene-d10	14.727	164	153205	20.00	ng	0.00
64) Phenanthrene-d10	17.471	188	432696	20.00	ng	# 0.00
76) Chrysene-d12	21.743	240	481607	20.00	ng	#-0.01
86) Perylene-d12	25.004	264	474673	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.661	112	346565	101.83	ng	0.00
7) Phenol-d6	7.254	99	491118	102.30	ng	0.00
23) Nitrobenzene-d5	9.263	82	364060	93.43	ng	0.00
42) 2,4,6-Tribromophenol	16.208	330	280553	138.88	ng	0.00
45) 2-Fluorobiphenyl	13.352	172	870084	80.40	ng	0.00
79) Terphenyl-d14	20.074	244	1824712	75.93	ng	0.00

Target Compounds	Qvalue

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

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