

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041321\
 Data File : BG048684.D
 Acq On : 13 Apr 2021 21:43
 Operator : CG/JU
 Sample : M1966-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EB 2021-04-06_GW

Quant Time: Apr 14 00:32:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 07 10:06:00 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.082	152	13867	20.00	ng	#-0.02
21) Naphthalene-d8	10.908	136	58456	20.00	ng	#-0.02
39) Acenaphthene-d10	14.733	164	43348	20.00	ng	-0.02
64) Phenanthrene-d10	17.489	188	116831	20.00	ng	-0.01
76) Chrysene-d12	21.784	240	124278	20.00	ng	-0.02
86) Perylene-d12	25.121	264	128954	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.632	112	61951	78.87	ng	0.00
7) Phenol-d6	7.242	99	50870	44.66	ng	0.00
23) Nitrobenzene-d5	9.251	82	111513	92.36	ng	-0.01
42) 2,4,6-Tribromophenol	16.231	330	79785	140.02	ng	0.00
45) 2-Fluorobiphenyl	13.358	172	247417	84.84	ng	-0.01
79) Terphenyl-d14	20.103	244	564683	83.09	ng	0.00
Target Compounds						
10) Phenol	7.265	94	2546	2.23	ng	# 80
50) Dimethylphthalate	14.181	163	14485	4.50	ng	# 98

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

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