

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041321\
 Data File : BG048683.D
 Acq On : 13 Apr 2021 21:03
 Operator : CG/JU
 Sample : M1967-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

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

Quant Time: Apr 14 00:32:43 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	8.081	152	15113	20.00	ng	#-0.02
21) Naphthalene-d8	10.913	136	63393	20.00	ng	#-0.01
39) Acenaphthene-d10	14.738	164	45480	20.00	ng	-0.01
64) Phenanthrene-d10	17.493	188	129568	20.00	ng	0.00
76) Chrysene-d12	21.788	240	131162	20.00	ng	#-0.01
86) Perylene-d12	25.126	264	133145	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.631	112	58073	67.84	ng	0.00
7) Phenol-d6	7.241	99	46891	37.77	ng	0.00
23) Nitrobenzene-d5	9.256	82	130224	99.46	ng	0.00
42) 2,4,6-Tribromophenol	16.230	330	95077	159.03	ng	0.00
45) 2-Fluorobiphenyl	13.357	172	291144	95.15	ng	-0.01
79) Terphenyl-d14	20.102	244	680274	94.85	ng	0.00

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

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