

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102921\
 Data File : BG050781.D
 Acq On : 29 Oct 2021 17:18
 Operator : CG/JU
 Sample : M4380-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WASTE-CHAR-(SUSPECT)

Quant Time: Oct 29 18:52:19 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 14 10:56:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.242	152	39855	20.000	ng	-0.02
21) Naphthalene-d8	11.068	136	175446	20.000	ng	#-0.02
39) Acenaphthene-d10	14.869	164	115222	20.000	ng	0.00
64) Phenanthrene-d10	17.613	188	261699	20.000	ng	# 0.00
76) Chrysene-d12	21.920	240	244600	20.000	ng	# 0.00
86) Perylene-d12	25.339	264	236665	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.786	112	287812	108.000	ng	0.00
7) Phenol-d6	7.390	99	338173	87.647	ng	-0.01
23) Nitrobenzene-d5	9.417	82	311004	73.380	ng	-0.01
42) 2,4,6-Tribromophenol	16.356	330	134955	122.796	ng	0.00
45) 2-Fluorobiphenyl	13.494	172	566753	74.029	ng	-0.01
79) Terphenyl-d14	20.204	244	946483	75.410	ng	0.00

Target Compounds	Qvalue

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

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