

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070221\
 Data File : BG049170.D
 Acq On : 2 Jul 2021 16:48
 Operator : CG/JU
 Sample : M2908-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-CROPSEY-SOLIDS

Quant Time: Jul 02 18:09:03 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG063021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 30 15:39:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.086	152	34830	20.00	ng	# 0.00
21) Naphthalene-d8	10.906	136	136790	20.00	ng	# 0.00
39) Acenaphthene-d10	14.731	164	100297	20.00	ng	0.00
64) Phenanthrene-d10	17.481	188	237356	20.00	ng	# 0.00
76) Chrysene-d12	21.764	240	236234	20.00	ng	-0.01
86) Perylene-d12	25.066	264	241425	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.642	112	271879	122.26	ng	0.00
7) Phenol-d6	7.240	99	325985	102.82	ng	0.00
23) Nitrobenzene-d5	9.255	82	279410	86.60	ng	0.00
42) 2,4,6-Tribromophenol	16.217	330	242094	139.24	ng	0.00
45) 2-Fluorobiphenyl	13.350	172	599231	80.61	ng	0.00
79) Terphenyl-d14	20.089	244	1163362	83.45	ng	0.00

Target Compounds	Qvalue

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

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