

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111721\
 Data File : BG051091.D
 Acq On : 17 Nov 2021 14:15
 Operator : CG/JU
 Sample : M4666-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CONCRETE

Quant Time: Nov 17 14:49:45 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 03:55:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.223	152	26512	20.000	ng	# 0.00
21) Naphthalene-d8	11.055	136	110581	20.000	ng	# 0.00
39) Acenaphthene-d10	14.851	164	77080	20.000	ng	0.00
64) Phenanthrene-d10	17.600	188	182369	20.000	ng	# 0.00
76) Chrysene-d12	21.901	240	156668	20.000	ng	# 0.00
86) Perylene-d12	25.315	264	160614	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.767	112	12979	7.323	ng	0.00
7) Phenol-d6	7.371	99	92100	36.818	ng	0.00
23) Nitrobenzene-d5	9.398	82	124994	50.036	ng	0.00
42) 2,4,6-Tribromophenol	16.343	330	1412	1.831	ng	0.00
45) 2-Fluorobiphenyl	13.476	172	255045	52.281	ng	0.00
79) Terphenyl-d14	20.186	244	434118	50.702	ng	0.00

Target Compounds	Qvalue

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

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