

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050624\
 Data File : BG061117.D
 Acq On : 6 May 2024 12:16
 Operator : MA/JU
 Sample : PB160699BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB160699BL

Quant Time: May 06 13:18:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG043024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 05:11:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.936	152	18304	20.000	ng	-0.02
21) Naphthalene-d8	10.733	136	69161	20.000	ng	#-0.02
39) Acenaphthene-d10	14.564	164	60869	20.000	ng	-0.02
64) Phenanthrene-d10	17.314	188	170438	20.000	ng	-0.02
76) Chrysene-d12	21.568	240	197451	20.000	ng	-0.02
86) Perylene-d12	24.699	264	222555	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	134549	149.496	ng	-0.02
7) Phenol-d6	7.114	99	177509	133.324	ng	-0.02
23) Nitrobenzene-d5	9.094	82	123583	88.580	ng	-0.02
42) 2,4,6-Tribromophenol	16.056	330	162596	133.641	ng	-0.02
45) 2-Fluorobiphenyl	13.189	172	357896	86.945	ng	-0.02
79) Terphenyl-d14	19.934	244	1056890	89.022	ng	-0.02

Target Compounds	Qvalue

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

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