

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102922\
 Data File : BG055369.D
 Acq On : 29 Oct 2022 17:27
 Operator : CG/JU
 Sample : N5318-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20221106-COMP-05

Quant Time: Oct 31 02:44:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 29 04:52:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.261	152	28132	20.000	ng	0.00
21) Naphthalene-d8	11.093	136	115486	20.000	ng	0.00
39) Acenaphthene-d10	14.871	164	78587	20.000	ng	0.00
64) Phenanthrene-d10	17.603	188	186634	20.000	ng	# 0.00
76) Chrysene-d12	21.881	240	198348	20.000	ng	0.00
86) Perylene-d12	25.271	264	214911	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.788	112	187590	116.758	ng	0.00
7) Phenol-d6	7.404	99	239213	102.770	ng	0.00
23) Nitrobenzene-d5	9.437	82	142194	62.873	ng	0.00
42) 2,4,6-Tribromophenol	16.352	330	115645	135.367	ng	0.00
45) 2-Fluorobiphenyl	13.502	172	333132	69.382	ng	0.00
79) Terphenyl-d14	20.177	244	630024	65.995	ng	0.00

Target Compounds Qvalue

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

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