

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112223\
 Data File : BG059847.D
 Acq On : 23 Nov 2023 11:17
 Operator : MA/JU
 Sample : 05449-07
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-SED-03

Quant Time: Nov 24 00:19:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 17 01:51:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.329	152	35036	20.000	ng	#-0.01
21) Naphthalene-d8	11.178	136	107037	20.000	ng	#-0.01
39) Acenaphthene-d10	14.962	164	90769	20.000	ng	0.00
64) Phenanthrene-d10	17.712	188	216073	20.000	ng	# 0.00
76) Chrysene-d12	22.042	240	169509	20.000	ng	# 0.00
86) Perylene-d12	25.626	264	202798	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.826	112	359417	172.665	ng	0.00
7) Phenol-d6	7.459	99	541552	175.560	ng	0.00
23) Nitrobenzene-d5	9.533	82	441052	157.372	ng	0.00
42) 2,4,6-Tribromophenol	16.449	330	237705	230.807	ng	0.00
45) 2-Fluorobiphenyl	13.587	172	714031	105.886	ng	0.00
79) Terphenyl-d14	20.291	244	1156835	100.748	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	9.081	107	7466	2.297	ng	Qvalue # 69

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

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