

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092322\
 Data File : BG055019.D
 Acq On : 23 Sep 2022 19:31
 Operator : CG/JU
 Sample : N4751-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PE-5D

Quant Time: Sep 24 02:29:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG091922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 19 17:41:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.059	152	41567	20.000	ng	0.00
21) Naphthalene-d8	10.885	136	165047	20.000	ng	# 0.00
39) Acenaphthene-d10	14.704	164	114690	20.000	ng	0.00
64) Phenanthrene-d10	17.448	188	262746	20.000	ng	0.00
76) Chrysene-d12	21.725	240	265706	20.000	ng	0.00
86) Perylene-d12	25.033	264	277058	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.597	112	235717	110.959	ng	0.00
7) Phenol-d6	7.218	99	306467	107.629	ng	0.00
23) Nitrobenzene-d5	9.234	82	177870	62.465	ng	0.00
42) 2,4,6-Tribromophenol	16.190	330	134022	99.749	ng	0.00
45) 2-Fluorobiphenyl	13.323	172	488570	67.135	ng	0.00
79) Terphenyl-d14	20.050	244	840772	66.191	ng	0.00

Target Compounds	Qvalue

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

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