

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081022\
 Data File : BG054590.D
 Acq On : 10 Aug 2022 16:07
 Operator : CG/JU
 Sample : N4117-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 080922-A

Quant Time: Aug 10 17:09:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 08 14:48:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.976	152	22722	20.000	ng	0.00
21) Naphthalene-d8	10.784	136	85554	20.000	ng	# 0.00
39) Acenaphthene-d10	14.615	164	62948	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	153047	20.000	ng	# 0.00
76) Chrysene-d12	21.631	240	169881	20.000	ng	# 0.00
86) Perylene-d12	24.844	264	179930	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.567	112	38321	35.212	ng	0.00
7) Phenol-d6	7.183	99	63625	42.667	ng	0.00
23) Nitrobenzene-d5	9.139	82	88517	56.343	ng	0.00
42) 2,4,6-Tribromophenol	16.113	330	35288	26.699	ng	0.00
45) 2-Fluorobiphenyl	13.234	172	257788	57.771	ng	0.00
79) Terphenyl-d14	19.968	244	510117	59.569	ng	0.00
Target Compounds						
25) Isophorone	9.703	82	8902	3.258	ng	# 92

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

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