

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050222\
 Data File : BG053397.D
 Acq On : 2 May 2022 13:44
 Operator : CG/JU
 Sample : N2616-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-20S

Quant Time: May 02 16:47:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 01:13:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.102	152	27044	20.000	ng	-0.01
21) Naphthalene-d8	10.915	136	100789	20.000	ng	# 0.00
39) Acenaphthene-d10	14.728	164	72183	20.000	ng	0.00
64) Phenanthrene-d10	17.471	188	187342	20.000	ng	-0.01
76) Chrysene-d12	21.748	240	213466	20.000	ng	-0.02
86) Perylene-d12	25.049	264	213638	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.670	112	224045	142.011	ng	0.00
7) Phenol-d6	7.267	99	288574	118.612	ng	0.00
23) Nitrobenzene-d5	9.265	82	241233	97.182	ng	-0.01
42) 2,4,6-Tribromophenol	16.214	330	172866	150.527	ng	-0.01
45) 2-Fluorobiphenyl	13.347	172	507450	96.938	ng	-0.01
79) Terphenyl-d14	20.074	244	1160093	91.784	ng	-0.01
Target Compounds						
77) Benzidine	19.692	184	76505	12.561	ng	Qvalue 98

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

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