

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091222\
 Data File : BG054911.D
 Acq On : 12 Sep 2022 18:12
 Operator : CG/JU
 Sample : N4555-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 257248-4

Quant Time: Sep 13 03:06:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 10 01:53:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.137	152	33479	20.000	ng	0.00
21) Naphthalene-d8	10.969	136	132914	20.000	ng	# 0.00
39) Acenaphthene-d10	14.776	164	97326	20.000	ng	0.00
64) Phenanthrene-d10	17.526	188	237033	20.000	ng	0.00
76) Chrysene-d12	21.815	240	258197	20.000	ng	0.00
86) Perylene-d12	25.194	264	267808	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.681	112	111134	57.805	ng	0.00
7) Phenol-d6	7.297	99	83915	31.916	ng	0.00
23) Nitrobenzene-d5	9.312	82	214314	84.647	ng	0.00
42) 2,4,6-Tribromophenol	16.269	330	182620	150.162	ng	0.00
45) 2-Fluorobiphenyl	13.396	172	583450	90.103	ng	0.00
79) Terphenyl-d14	20.123	244	1086149	83.350	ng	0.00

Target Compounds Qvalue

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

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