

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110222\
 Data File : BG055427.D
 Acq On : 2 Nov 2022 23:20
 Operator : CG/JU
 Sample : N5396-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-G

Quant Time: Nov 03 06:13:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 02 05:05:12 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.240	152	31943	20.000	ng	0.00
21) Naphthalene-d8	11.072	136	135049	20.000	ng	0.00
39) Acenaphthene-d10	14.856	164	91864	20.000	ng	0.00
64) Phenanthrene-d10	17.588	188	191943	20.000	ng	0.00
76) Chrysene-d12	21.866	240	186484	20.000	ng	0.00
86) Perylene-d12	25.238	264	200792	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.767	112	229421	128.981	ng	0.00
7) Phenol-d6	7.389	99	273860	111.026	ng	0.00
23) Nitrobenzene-d5	9.416	82	206884	81.643	ng	0.00
42) 2,4,6-Tribromophenol	16.337	330	163101	130.164	ng	0.00
45) 2-Fluorobiphenyl	13.481	172	534342	86.241	ng	0.00
79) Terphenyl-d14	20.168	244	844542	88.063	ng	0.00

Target Compounds	Qvalue

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

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