

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110222\
 Data File : BG055426.D
 Acq On : 2 Nov 2022 22:38
 Operator : CG/JU
 Sample : N5396-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-F

Quant Time: Nov 03 06:12:51 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	8.241	152	32540	20.000	ng	0.00
21) Naphthalene-d8	11.068	136	140399	20.000	ng	0.00
39) Acenaphthene-d10	14.857	164	94251	20.000	ng	0.00
64) Phenanthrene-d10	17.589	188	197304	20.000	ng	0.00
76) Chrysene-d12	21.861	240	189350	20.000	ng	-0.01
86) Perylene-d12	25.233	264	205377	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.768	112	235189	129.798	ng	0.00
7) Phenol-d6	7.390	99	279071	111.063	ng	0.00
23) Nitrobenzene-d5	9.417	82	213413	81.011	ng	0.00
42) 2,4,6-Tribromophenol	16.338	330	165490	128.725	ng	0.00
45) 2-Fluorobiphenyl	13.482	172	548847	86.339	ng	0.00
79) Terphenyl-d14	20.163	244	859313	88.247	ng	0.00

Target Compounds	Qvalue

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

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