

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110222\
 Data File : BG055431.D
 Acq On : 3 Nov 2022 2:06
 Operator : CG/JU
 Sample : N5396-14
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-I

Quant Time: Nov 03 06:14:55 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	8.242	152	34509	20.000	ng	0.00
21) Naphthalene-d8	11.068	136	151435	20.000	ng	0.00
39) Acenaphthene-d10	14.857	164	101727	20.000	ng	0.00
64) Phenanthrene-d10	17.589	188	216587	20.000	ng	0.00
76) Chrysene-d12	21.861	240	208694	20.000	ng	-0.01
86) Perylene-d12	25.233	264	225896	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.768	112	244408	127.190	ng	0.00
7) Phenol-d6	7.390	99	293357	110.087	ng	0.00
23) Nitrobenzene-d5	9.417	82	221832	78.070	ng	0.00
42) 2,4,6-Tribromophenol	16.338	330	177677	128.048	ng	0.00
45) 2-Fluorobiphenyl	13.482	172	577469	84.165	ng	0.00
79) Terphenyl-d14	20.163	244	920463	85.765	ng	0.00
Target Compounds						
70) Pentachlorophenol	17.278	266	6868	8.506	ng	Qvalue 89

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

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