

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040622\
 Data File : BG053056.D
 Acq On : 6 Apr 2022 15:50
 Operator : CG/JU
 Sample : N2269-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LP-2

Quant Time: Apr 06 16:30:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 06 14:35:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.128	152	32374	20.000	ng	0.00
21) Naphthalene-d8	10.936	136	130101	20.000	ng	0.00
39) Acenaphthene-d10	14.748	164	94288	20.000	ng	0.00
64) Phenanthrene-d10	17.492	188	224777	20.000	ng	0.00
76) Chrysene-d12	21.774	240	223349	20.000	ng	0.00
86) Perylene-d12	25.082	264	229922	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.696	112	147118	79.140	ng	0.01
7) Phenol-d6	7.288	99	174680	62.328	ng	0.00
23) Nitrobenzene-d5	9.285	82	278427	107.411	ng	0.00
42) 2,4,6-Tribromophenol	16.235	330	219640	173.357	ng	0.00
45) 2-Fluorobiphenyl	13.374	172	606463	94.863	ng	0.00
79) Terphenyl-d14	20.094	244	1164673	92.760	ng	0.00

Target Compounds	Qvalue

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

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