

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020723\
 Data File : BG056568.D
 Acq On : 7 Feb 2023 15:18
 Operator : CG/JU
 Sample : 01462-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COMP-2

Quant Time: Feb 08 03:05:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG012723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 08 03:02:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.268	152	36995	20.000	ng	# 0.00
21) Naphthalene-d8	11.100	136	140085	20.000	ng	0.00
39) Acenaphthene-d10	14.889	164	106975	20.000	ng	0.00
64) Phenanthrene-d10	17.627	188	261051	20.000	ng	0.00
76) Chrysene-d12	21.916	240	239399	20.000	ng	0.00
86) Perylene-d12	25.336	264	261479	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.794	112	282003	140.226	ng	0.00
7) Phenol-d6	7.422	99	383654	128.755	ng	0.00
23) Nitrobenzene-d5	9.443	82	202021	81.891	ng	0.00
42) 2,4,6-Tribromophenol	16.376	330	183780	129.225	ng	0.00
45) 2-Fluorobiphenyl	13.515	172	668635	86.657	ng	0.00
79) Terphenyl-d14	20.207	244	1184612	86.910	ng	0.00
Target Compounds						
75) Fluoranthene	19.660	202	53669	3.112	ng	Qvalue 98
78) Pyrene	20.025	202	52967	3.111	ng	98
88) Benzo(b)fluoranthene	24.249	252	38232	2.356	ng	# 93

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

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