

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080122\
 Data File : BG054470.D
 Acq On : 1 Aug 2022 18:44
 Operator : CG/JU
 Sample : N3896-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-59

Quant Time: Aug 02 05:46:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 04:30:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.982	152	17975	20.000	ng	0.00
21) Naphthalene-d8	10.790	136	70879	20.000	ng	# 0.00
39) Acenaphthene-d10	14.621	164	59032	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	139267	20.000	ng	# 0.00
76) Chrysene-d12	21.630	240	153454	20.000	ng	# 0.00
86) Perylene-d12	24.832	264	163139	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.567	112	88955	103.326	ng	0.00
7) Phenol-d6	7.177	99	122890	104.173	ng	0.00
23) Nitrobenzene-d5	9.145	82	84812	65.162	ng	0.00
42) 2,4,6-Tribromophenol	16.113	330	123950	100.004	ng	0.00
45) 2-Fluorobiphenyl	13.240	172	279714	66.844	ng	0.00
79) Terphenyl-d14	19.973	244	554918	71.738	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.421	202	38046	4.024	ng	96
78) Pyrene	19.785	202	33766	3.835	ng	99
88) Benzo(b)fluoranthene	23.816	252	29834	3.092	ng	# 97
90) Benzo(a)pyrene	24.686	252	18453	2.240	ng	# 87

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

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