

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080122\
 Data File : BG054466.D
 Acq On : 1 Aug 2022 15:58
 Operator : CG/JU
 Sample : N3896-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-66

Quant Time: Aug 02 05:45:56 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	15487	20.000	ng	0.00
21) Naphthalene-d8	10.790	136	61935	20.000	ng	# 0.00
39) Acenaphthene-d10	14.621	164	52023	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	133675	20.000	ng	# 0.00
76) Chrysene-d12	21.630	240	147043	20.000	ng	# 0.00
86) Perylene-d12	24.833	264	155474	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.561	112	77892	105.010	ng	0.00
7) Phenol-d6	7.177	99	107986	106.245	ng	0.00
23) Nitrobenzene-d5	9.145	82	74999	65.944	ng	0.00
42) 2,4,6-Tribromophenol	16.113	330	117516	107.587	ng	0.00
45) 2-Fluorobiphenyl	13.240	172	247565	67.132	ng	0.00
79) Terphenyl-d14	19.974	244	560919	75.675	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.421	202	46614	5.137	ng	95
78) Pyrene	19.786	202	42377	5.023	ng	97
81) Benzo(a)anthracene	21.613	228	26596	2.849	ng	96
83) Chrysene	21.677	228	27945	3.169	ng	98
88) Benzo(b)fluoranthene	23.822	252	44174	4.804	ng	# 94
90) Benzo(a)pyrene	24.686	252	24970	3.180	ng	96

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

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