

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080122\
 Data File : BG054468.D
 Acq On : 1 Aug 2022 17:21
 Operator : CG/JU
 Sample : N3894-19
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-38

Quant Time: Aug 02 05:46:12 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	7.984	152	17323	20.000	ng	0.00
21) Naphthalene-d8	10.792	136	67694	20.000	ng	# 0.00
39) Acenaphthene-d10	14.617	164	56096	20.000	ng	0.00
64) Phenanthrene-d10	17.367	188	142050	20.000	ng	# 0.00
76) Chrysene-d12	21.633	240	157580	20.000	ng	# 0.00
86) Perylene-d12	24.835	264	167867	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	93258	112.401	ng	0.00
7) Phenol-d6	7.173	99	129601	113.997	ng	0.00
23) Nitrobenzene-d5	9.147	82	86263	69.395	ng	0.00
42) 2,4,6-Tribromophenol	16.115	330	134284	114.012	ng	0.00
45) 2-Fluorobiphenyl	13.242	172	282124	70.948	ng	0.00
79) Terphenyl-d14	19.976	244	600806	75.637	ng	0.00
Target Compounds						
						Qvalue
78) Pyrene	19.782	202	38517	4.260	ng	97
88) Benzo(b)fluoranthene	23.824	252	24149	2.433	ng	# 97
90) Benzo(a)pyrene	24.682	252	20013	2.361	ng	# 94

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

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