

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080422\
 Data File : BG054516.D
 Acq On : 4 Aug 2022 12:49
 Operator : CG/JU
 Sample : N3931-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-119

Quant Time: Aug 04 13:44:59 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	7.975	152	18170	20.000	ng	-0.01
21) Naphthalene-d8	10.790	136	68741	20.000	ng	# 0.00
39) Acenaphthene-d10	14.615	164	54684	20.000	ng	0.00
64) Phenanthrene-d10	17.364	188	138533	20.000	ng	# 0.00
76) Chrysene-d12	21.630	240	158307	20.000	ng	# 0.00
86) Perylene-d12	24.838	264	162413	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	103950	119.447	ng	0.00
7) Phenol-d6	7.182	99	139325	116.837	ng	0.01
23) Nitrobenzene-d5	9.145	82	91260	72.297	ng	0.00
42) 2,4,6-Tribromophenol	16.113	330	125338	109.164	ng	0.00
45) 2-Fluorobiphenyl	13.240	172	275187	70.991	ng	0.00
79) Terphenyl-d14	19.973	244	591402	74.111	ng	0.00
Target Compounds						
88) Benzo(b)fluoranthene	23.822	252	22461	2.338	ng	Qvalue 99

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

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