

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080522\
 Data File : BG054536.D
 Acq On : 5 Aug 2022 15:26
 Operator : CG/JU
 Sample : N3960-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-155

Quant Time: Aug 06 05:39:45 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	7.976	152	16166	20.000	ng	-0.01
21) Naphthalene-d8	10.785	136	68345	20.000	ng	#-0.01
39) Acenaphthene-d10	14.616	164	55771	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	137178	20.000	ng	# 0.00
76) Chrysene-d12	21.631	240	145314	20.000	ng	# 0.00
86) Perylene-d12	24.839	264	156956	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.561	112	86191	111.318	ng	0.00
7) Phenol-d6	7.177	99	122898	115.838	ng	0.00
23) Nitrobenzene-d5	9.140	82	82458	65.702	ng	-0.01
42) 2,4,6-Tribromophenol	16.114	330	114527	97.804	ng	0.00
45) 2-Fluorobiphenyl	13.235	172	264823	66.985	ng	-0.01
79) Terphenyl-d14	19.974	244	528586	72.162	ng	0.00
Target Compounds						
						Qvalue
68) Hexachlorobenzene	16.678	284	10355	4.700	ng	# 93
75) Fluoranthene	19.416	202	33236	3.569	ng	97
78) Pyrene	19.780	202	36639	4.394	ng	96
83) Chrysene	21.678	228	22535	2.586	ng	97
88) Benzo(b)fluoranthene	23.822	252	32255	3.475	ng	# 95
90) Benzo(a)pyrene	24.686	252	21655	2.732	ng	# 90

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

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