

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080422\
 Data File : BG054517.D
 Acq On : 4 Aug 2022 13:30
 Operator : CG/JU
 Sample : N3931-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-119D

Quant Time: Aug 04 14:45:28 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	19234	20.000	ng	-0.01
21) Naphthalene-d8	10.790	136	71586	20.000	ng	# 0.00
39) Acenaphthene-d10	14.615	164	55107	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	137775	20.000	ng	# 0.00
76) Chrysene-d12	21.630	240	152793	20.000	ng	# 0.00
86) Perylene-d12	24.832	264	160211	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.567	112	93987	102.025	ng	0.00
7) Phenol-d6	7.177	99	124586	98.698	ng	0.00
23) Nitrobenzene-d5	9.145	82	82787	62.978	ng	0.00
42) 2,4,6-Tribromophenol	16.113	330	108697	93.944	ng	0.00
45) 2-Fluorobiphenyl	13.234	172	248204	63.538	ng	-0.01
79) Terphenyl-d14	19.973	244	508075	65.966	ng	0.00
Target Compounds						
75) Fluoranthene	19.415	202	24797	2.651	ng	Qvalue 96
78) Pyrene	19.780	202	24752	2.823	ng	97
88) Benzo(b)fluoranthene	23.816	252	28493	3.007	ng	96
90) Benzo(a)pyrene	24.680	252	18605	2.299	ng	# 96

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

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