

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080422\
 Data File : BG054519.D
 Acq On : 4 Aug 2022 14:57
 Operator : CG/JU
 Sample : N4037-04 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 080322-C

Quant Time: Aug 04 15:32:27 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	24907	20.000	ng	0.00
21) Naphthalene-d8	10.790	136	92751	20.000	ng	# 0.00
39) Acenaphthene-d10	14.621	164	67112	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	155041	20.000	ng	# 0.00
76) Chrysene-d12	21.636	240	172510	20.000	ng	# 0.00
86) Perylene-d12	24.856	264	184124	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.573	112	6189	5.188	ng	0.01
7) Phenol-d6	7.189	99	9850	6.026	ng	0.02
23) Nitrobenzene-d5	9.151	82	6956	4.084	ng	0.00
42) 2,4,6-Tribromophenol	16.119	330	3730	2.647	ng	0.00
45) 2-Fluorobiphenyl	13.234	172	22850	4.803	ng	-0.01
79) Terphenyl-d14	19.974	244	47323	5.442	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.412	178	21496	2.784	ng	98
75) Fluoranthene	19.421	202	29431	2.796	ng	91
78) Pyrene	19.786	202	27298	2.758	ng	96

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

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