

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG103122\
 Data File : BG055390.D
 Acq On : 31 Oct 2022 22:42
 Operator : CG/JU
 Sample : N5371-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-2-WC

Quant Time: Nov 01 06:51:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 29 04:52:38 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.258	152	35603	20.000	ng	0.00
21) Naphthalene-d8	11.090	136	141905	20.000	ng	# 0.00
39) Acenaphthene-d10	14.874	164	96614	20.000	ng	0.00
64) Phenanthrene-d10	17.606	188	214636	20.000	ng	# 0.00
76) Chrysene-d12	21.884	240	209865	20.000	ng	# 0.00
86) Perylene-d12	25.268	264	225189	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.791	112	246470	121.215	ng	0.00
7) Phenol-d6	7.406	99	290812	98.720	ng	0.00
23) Nitrobenzene-d5	9.433	82	220041	79.180	ng	0.00
42) 2,4,6-Tribromophenol	16.355	330	171399	163.194	ng	0.00
45) 2-Fluorobiphenyl	13.499	172	527059	89.290	ng	-0.01
79) Terphenyl-d14	20.180	244	911676	90.257	ng	0.00
Target Compounds						
31) Naphthalene	11.137	128	75771	10.009	ng	Qvalue 99

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

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