

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112922\
 Data File : BG055837.D
 Acq On : 29 Nov 2022 20:50
 Operator : CG/JU
 Sample : N5787-03 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-1-STOCK

Quant Time: Nov 30 02:06:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 30 02:01:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.224	152	56972	20.000	ng	0.00
21) Naphthalene-d8	11.050	136	236602	20.000	ng	0.00
39) Acenaphthene-d10	14.846	164	163105	20.000	ng	0.00
64) Phenanthrene-d10	17.590	188	353073	20.000	ng	0.00
76) Chrysene-d12	21.879	240	331411	20.000	ng	0.00
86) Perylene-d12	25.269	264	354790	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.757	112	62568	19.851	ng	0.00
7) Phenol-d6	7.372	99	83558	19.915	ng	0.00
23) Nitrobenzene-d5	9.394	82	51553	11.516	ng	0.00
42) 2,4,6-Tribromophenol	16.333	330	39903	17.374	ng	0.00
45) 2-Fluorobiphenyl	13.471	172	134911	12.392	ng	0.00
79) Terphenyl-d14	20.169	244	223534	13.491	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.629	202	84070	3.627	ng	98
78) Pyrene	19.987	202	81254	3.622	ng	99
81) Benzo(a)anthracene	21.855	228	56619	2.480	ng	98
83) Chrysene	21.926	228	54451	2.506	ng	97
88) Benzo(b)fluoranthene	24.188	252	75742	3.278	ng	97
90) Benzo(a)pyrene	25.105	252	54085	2.728	ng	# 93

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

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