

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112423\
 Data File : BG059885.D
 Acq On : 25 Nov 2023 10:35
 Operator : MA/JU
 Sample : 05366-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P-13(10-15)

Quant Time: Nov 27 03:19:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 24 15:11:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.312	152	22507	20.000	ng	0.00
21) Naphthalene-d8	11.167	136	119268	20.000	ng	# 0.00
39) Acenaphthene-d10	14.951	164	82940	20.000	ng	# 0.00
64) Phenanthrene-d10	17.695	188	160895	20.000	ng	-0.01
76) Chrysene-d12	22.019	240	128980	20.000	ng	#-0.01
86) Perylene-d12	25.580	264	164508	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.809	112	124281	92.941	ng	0.00
7) Phenol-d6	7.436	99	142349	71.835	ng	0.00
23) Nitrobenzene-d5	9.516	82	164695	52.739	ng	0.00
42) 2,4,6-Tribromophenol	16.432	330	75786	80.533	ng	0.00
45) 2-Fluorobiphenyl	13.570	172	322400	52.323	ng	0.00
79) Terphenyl-d14	20.274	244	508360	58.185	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	11.214	128	13292	2.216	ng	# 92
71) Phenanthrene	17.742	178	28441	3.477	ng	# 94
78) Pyrene	20.092	202	23922	2.726	ng	94

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

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