

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
 Data File : BG059758.D
 Acq On : 18 Nov 2023 10:31
 Operator : MA/JU
 Sample : 05374-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RB21197

Quant Time: Nov 20 04:41:00 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 17 01:51:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.331	152	17434	20.000	ng	-0.01
21) Naphthalene-d8	11.192	136	78355	20.000	ng	# 0.00
39) Acenaphthene-d10	14.970	164	63426	20.000	ng	0.00
64) Phenanthrene-d10	17.720	188	191501	20.000	ng	# 0.00
76) Chrysene-d12	22.050	240	168624	20.000	ng	# 0.00
86) Perylene-d12	25.634	264	216415	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.834	112	86921	83.917	ng	0.00
7) Phenol-d6	7.461	99	139245	90.716	ng	0.00
23) Nitrobenzene-d5	9.541	82	124534	60.701	ng	0.00
42) 2,4,6-Tribromophenol	16.451	330	96845	134.573	ng	0.00
45) 2-Fluorobiphenyl	13.589	172	386583	82.042	ng	0.00
79) Terphenyl-d14	20.293	244	646827	56.628	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.761	178	62589	6.428	ng	# 92
75) Fluoranthene	19.753	202	61490	5.266	ng	90
78) Pyrene	20.117	202	45632	3.977	ng	94

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

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