

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061821\
 Data File : BG049021.D
 Acq On : 18 Jun 2021 15:49
 Operator : CG/JU
 Sample : M2687-12
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 17-B6(104-108)

Quant Time: Jun 18 16:24:32 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 16 12:04:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.115	152	29286	20.000	ng	# 0.00
21) Naphthalene-d8	10.935	136	116238	20.000	ng	# 0.00
39) Acenaphthene-d10	14.754	164	77398	20.000	ng	0.00
64) Phenanthrene-d10	17.504	188	189235	20.000	ng	# 0.00
76) Chrysene-d12	21.793	240	245946	20.000	ng	0.00
86) Perylene-d12	25.118	264	253027	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.664	112	265639	141.199	ng	0.00
7) Phenol-d6	7.263	99	319136	113.180	ng	0.00
23) Nitrobenzene-d5	9.278	82	262749	97.470	ng	0.00
42) 2,4,6-Tribromophenol	16.240	330	187374	163.403	ng	0.00
45) 2-Fluorobiphenyl	13.379	172	541694	98.870	ng	0.00
79) Terphenyl-d14	20.112	244	1219936	90.841	ng	0.00

Target Compounds	Qvalue

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

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