

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061821\
 Data File : BG049022.D
 Acq On : 18 Jun 2021 16:31
 Operator : CG/JU
 Sample : M2687-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 17-B6(108-112)

Quant Time: Jun 18 17:06:25 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	8.114	152	28229	20.000	ng	0.00
21) Naphthalene-d8	10.934	136	110679	20.000	ng	# 0.00
39) Acenaphthene-d10	14.753	164	77220	20.000	ng	0.00
64) Phenanthrene-d10	17.503	188	192038	20.000	ng	# 0.00
76) Chrysene-d12	21.792	240	240833	20.000	ng	0.00
86) Perylene-d12	25.111	264	251582	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.664	112	250805	138.306	ng	0.00
7) Phenol-d6	7.262	99	312000	114.792	ng	0.00
23) Nitrobenzene-d5	9.277	82	253380	98.715	ng	0.00
42) 2,4,6-Tribromophenol	16.239	330	190045	166.114	ng	0.00
45) 2-Fluorobiphenyl	13.378	172	529257	96.822	ng	0.00
79) Terphenyl-d14	20.111	244	1247252	94.847	ng	0.00

Target Compounds	Qvalue

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

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