

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061821\
 Data File : BG049018.D
 Acq On : 18 Jun 2021 13:45
 Operator : CG/JU
 Sample : M2687-11
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 17-B6(100-104)

Quant Time: Jun 18 14:21:50 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	8.109	152	30035	20.000	ng	0.00
21) Naphthalene-d8	10.930	136	108379	20.000	ng	#-0.01
39) Acenaphthene-d10	14.754	164	74041	20.000	ng	0.00
64) Phenanthrene-d10	17.498	188	172787	20.000	ng	#-0.01
76) Chrysene-d12	21.787	240	230340	20.000	ng	-0.01
86) Perylene-d12	25.113	264	243728	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.665	112	242595	125.734	ng	0.00
7) Phenol-d6	7.269	99	279581	96.679	ng	0.00
23) Nitrobenzene-d5	9.279	82	245553	97.696	ng	0.00
42) 2,4,6-Tribromophenol	16.241	330	164436	149.901	ng	0.00
45) 2-Fluorobiphenyl	13.380	172	492966	94.055	ng	0.00
79) Terphenyl-d14	20.107	244	1048219	83.343	ng	0.00

Target Compounds	Qvalue

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

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