

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022222\
 Data File : BN018681.D
 Acq On : 22 Feb 2022 17:26
 Operator : CG/JU
 Sample : PB142611BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB142611BL

Quant Time: Feb 23 02:17:09 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 22 16:23:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.919	152	6693	0.400	ng	0.00
7) Naphthalene-d8	10.722	136	15815	0.400	ng	# 0.00
13) Acenaphthene-d10	14.549	164	7313	0.400	ng	0.00
19) Phenanthrene-d10	17.287	188	13300	0.400	ng	# 0.00
29) Chrysene-d12	21.467	240	10165	0.400	ng	# 0.00
35) Perylene-d12	23.861	264	9561	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.471	112	3948	0.258	ng	0.00
5) Phenol-d6	7.074	99	3837	0.218	ng	0.00
8) Nitrobenzene-d5	9.078	82	2966	0.243	ng	0.01
11) 2-Methylnaphthalene-d10	12.314	152	6056	0.254	ng	0.00
14) 2,4,6-Tribromophenol	16.036	330	432	0.151	ng	0.01
15) 2-Fluorobiphenyl	13.180	172	9339	0.269	ng	0.00
27) Fluoranthene-d10	19.316	212	9485	0.263	ng	0.00
31) Terphenyl-d14	19.906	244	6390	0.290	ng	0.00

Target Compounds Qvalue

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

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