

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112320\
 Data File : BN012696.D
 Acq On : 23 Nov 2020 14:41
 Operator : CG/JU
 Sample : PB133106BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB133106BL

Quant Time: Nov 23 16:21:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN110420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 12:44:50 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.441	152	504	0.40	ng	# 0.00
7) Naphthalene-d8	10.212	136	1941	0.40	ng	# 0.01
13) Acenaphthene-d10	14.090	164	1120	0.40	ng	0.00
19) Phenanthrene-d10	16.859	188	2271	0.40	ng	# 0.01
29) Chrysene-d12	21.067	240	2417	0.40	ng	# 0.00
36) Perylene-d12	23.183	264	2965	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.057	112	677	0.38	ng	0.00
5) Phenol-d6	6.636	99	659	0.32	ng	0.00
8) Nitrobenzene-d5	8.603	82	1149	0.50	ng	0.01
11) 2-Methylnaphthalene-d10	11.811	152	1093	0.35	ng	0.00
14) 2,4,6-Tribromophenol	15.613	330	282	0.35	ng	0.01
15) 2-Fluorobiphenyl	12.707	172	2240	0.51	ng	0.00
27) Fluoranthene-d10	18.893	212	2628	0.37	ng	0.00
31) Terphenyl-d14	19.509	244	3742	0.65	ng	0.00

Target Compounds Qvalue

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

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