

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070822\
 Data File : BN020688.D
 Acq On : 08 Jul 2022 11:54
 Operator : CG/JU
 Sample : PB146079BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB146079BL

Quant Time: Jul 09 00:09:05 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN062622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 02 03:10:12 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.825	152	4330	0.400	ng	0.00
7) Naphthalene-d8	10.626	136	12659	0.400	ng	# 0.00
13) Acenaphthene-d10	14.474	164	6865	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	12093	0.400	ng	# 0.01
29) Chrysene-d12	21.440	240	7408	0.400	ng	0.00
35) Perylene-d12	23.790	264	6142	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.413	112	3635	0.302	ng	0.00
5) Phenol-d6	7.031	99	3406	0.237	ng	0.00
8) Nitrobenzene-d5	9.014	82	2584	0.252	ng	0.02
11) 2-Methylnaphthalene-d10	12.231	152	6721	0.310	ng	0.00
14) 2,4,6-Tribromophenol	15.984	330	424	0.164	ng	0.01
15) 2-Fluorobiphenyl	13.105	172	8926	0.315	ng	0.00
27) Fluoranthene-d10	19.270	212	10181	0.283	ng	0.00
31) Terphenyl-d14	19.872	244	7288	0.412	ng	0.00

Target Compounds	Qvalue

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

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