

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070622\
 Data File : BN020669.D
 Acq On : 06 Jul 2022 16:16
 Operator : CG/JU
 Sample : PB146049BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB146049BL

Quant Time: Jul 06 23:48:08 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	7.818	152	4631	0.400	ng	0.00
7) Naphthalene-d8	10.616	136	13649	0.400	ng	#-0.01
13) Acenaphthene-d10	14.474	164	7525	0.400	ng	0.00
19) Phenanthrene-d10	17.225	188	14928	0.400	ng	# 0.00
29) Chrysene-d12	21.440	240	10486	0.400	ng	0.00
35) Perylene-d12	23.793	264	8881	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.399	112	3963	0.308	ng	-0.01
5) Phenol-d6	7.017	99	3860	0.251	ng	-0.01
8) Nitrobenzene-d5	8.993	82	2880	0.260	ng	0.00
11) 2-Methylnaphthalene-d10	12.219	152	7251	0.310	ng	0.00
14) 2,4,6-Tribromophenol	15.984	330	525	0.185	ng	0.01
15) 2-Fluorobiphenyl	13.095	172	10144	0.327	ng	-0.01
27) Fluoranthene-d10	19.270	212	12930	0.292	ng	0.00
31) Terphenyl-d14	19.872	244	9350	0.373	ng	0.00

Target Compounds Qvalue

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

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