

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071222\
 Data File : BN020713.D
 Acq On : 12 Jul 2022 12:47
 Operator : CG/JU
 Sample : PB146111BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB146111BL

Quant Time: Jul 13 02:41:16 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	7.826	152	4304	0.400	ng	0.00
7) Naphthalene-d8	10.626	136	12285	0.400	ng	# 0.00
13) Acenaphthene-d10	14.474	164	6631	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	12280	0.400	ng	# 0.01
29) Chrysene-d12	21.440	240	8210	0.400	ng	# 0.00
35) Perylene-d12	23.793	264	6662	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.406	112	3663	0.307	ng	0.00
5) Phenol-d6	7.038	99	3377	0.237	ng	0.00
8) Nitrobenzene-d5	9.014	82	2582	0.259	ng	0.02
11) 2-Methylnaphthalene-d10	12.231	152	6834	0.325	ng	0.00
14) 2,4,6-Tribromophenol	15.985	330	449	0.180	ng	0.01
15) 2-Fluorobiphenyl	13.105	172	8849	0.323	ng	0.00
27) Fluoranthene-d10	19.270	212	11592	0.318	ng	0.00
31) Terphenyl-d14	19.872	244	8196	0.418	ng	0.00

Target Compounds Qvalue

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

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