

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070822\
 Data File : BN020694.D
 Acq On : 08 Jul 2022 16:29
 Operator : CG/JU
 Sample : PB146111BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB146111BL

Quant Time: Jul 09 00:26:13 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	5023	0.400	ng	0.00
7) Naphthalene-d8	10.626	136	14830	0.400	ng	# 0.00
13) Acenaphthene-d10	14.474	164	8075	0.400	ng	0.00
19) Phenanthrene-d10	17.226	188	13134	0.400	ng	# 0.00
29) Chrysene-d12	21.440	240	6949	0.400	ng	0.00
35) Perylene-d12	23.793	264	6119	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.413	112	4066	0.292	ng	0.00
5) Phenol-d6	7.038	99	3967	0.238	ng	0.00
8) Nitrobenzene-d5	9.014	82	3090	0.257	ng	0.02
11) 2-Methylnaphthalene-d10	12.231	152	7955	0.313	ng	0.00
14) 2,4,6-Tribromophenol	15.985	330	481	0.158	ng	0.01
15) 2-Fluorobiphenyl	13.105	172	10753	0.323	ng	0.00
27) Fluoranthene-d10	19.270	212	10171	0.261	ng	0.00
31) Terphenyl-d14	19.872	244	7109	0.428	ng	0.00

Target Compounds Qvalue

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

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