

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031222\
 Data File : BN018931.D
 Acq On : 11 Mar 2022 19:18
 Operator : CG/JU
 Sample : PB143229BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143229BL

Quant Time: Mar 11 23:34:36 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN031222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 11 23:19:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.898	152	5522	0.400	ng	0.00
7) Naphthalene-d8	10.701	136	13652	0.400	ng	# 0.00
13) Acenaphthene-d10	14.527	164	7444	0.400	ng	0.00
19) Phenanthrene-d10	17.267	188	16239	0.400	ng	0.00
29) Chrysene-d12	21.449	240	10396	0.400	ng	# 0.00
35) Perylene-d12	23.837	264	8158	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.457	112	3983	0.287	ng	0.00
5) Phenol-d6	7.060	99	3919	0.242	ng	0.00
8) Nitrobenzene-d5	9.057	82	2783	0.246	ng	0.01
11) 2-Methylnaphthalene-d10	12.291	152	6095	0.271	ng	0.00
14) 2,4,6-Tribromophenol	16.026	330	671	0.173	ng	0.01
15) 2-Fluorobiphenyl	13.159	172	9196	0.284	ng	0.00
27) Fluoranthene-d10	19.299	212	12696	0.276	ng	0.00
31) Terphenyl-d14	19.889	244	7795	0.337	ng	0.00

Target Compounds Qvalue

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

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