

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031722\
 Data File : BN018993.D
 Acq On : 17 Mar 2022 14:15
 Operator : CG/JU
 Sample : PB143270BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143270BL

Quant Time: Mar 17 15:58:54 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 13:18:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.854	152	4958	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	12658	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	7553	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	13404	0.400	ng	0.00
29) Chrysene-d12	21.431	240	9850	0.400	ng	# 0.00
35) Perylene-d12	23.805	264	7118	0.400	ng	# 0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.406	112	3512	0.316	ng	0.00
5) Phenol-d6	7.009	99	3258	0.280	ng	0.00
8) Nitrobenzene-d5	9.003	82	2619	0.287	ng	-0.01
11) 2-Methylnaphthalene-d10	12.250	152	5694	0.271	ng	0.00
14) 2,4,6-Tribromophenol	15.984	330	553	0.152	ng	0.00
15) 2-Fluorobiphenyl	13.116	172	8870	0.293	ng	-0.01
27) Fluoranthene-d10	19.270	212	12592	0.317	ng	0.00
31) Terphenyl-d14	19.868	244	8825	0.365	ng	0.00

Target Compounds Qvalue

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

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