

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031822\
 Data File : BN019010.D
 Acq On : 18 Mar 2022 10:54
 Operator : CG/JU
 Sample : PB143277BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143277BL

Quant Time: Mar 18 15:38:01 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	6990	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	17234	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	10070	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	17291	0.400	ng	0.00
29) Chrysene-d12	21.422	240	13082	0.400	ng	# 0.00
35) Perylene-d12	23.793	264	10008	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.413	112	5339	0.341	ng	0.00
5) Phenol-d6	7.016	99	4528	0.276	ng	0.00
8) Nitrobenzene-d5	9.014	82	3627	0.292	ng	0.00
11) 2-Methylnaphthalene-d10	12.253	152	8659	0.303	ng	0.00
14) 2,4,6-Tribromophenol	15.995	330	668	0.137	ng	0.01
15) 2-Fluorobiphenyl	13.127	172	12813	0.317	ng	0.00
27) Fluoranthene-d10	19.265	212	19248	0.375	ng	0.00
31) Terphenyl-d14	19.864	244	11936	0.371	ng	0.00

Target Compounds Qvalue

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

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