

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031822\
 Data File : BN019015.D
 Acq On : 18 Mar 2022 14:50
 Operator : CG/JU
 Sample : PB143270BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143270BL

Quant Time: Mar 18 16:35:52 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 16:26:38 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.854	152	5659	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	14072	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	7852	0.400	ng	0.00
19) Phenanthrene-d10	17.246	188	12469	0.400	ng	0.00
29) Chrysene-d12	21.431	240	8948	0.400	ng	# 0.00
35) Perylene-d12	23.796	264	5904	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.413	112	4065	0.321	ng	0.00
5) Phenol-d6	7.016	99	3082	0.232	ng	0.00
8) Nitrobenzene-d5	9.024	82	2728	0.269	ng	0.01
11) 2-Methylnaphthalene-d10	12.257	152	7074	0.303	ng	0.00
14) 2,4,6-Tribromophenol	15.994	330	435	0.115	ng	0.00
15) 2-Fluorobiphenyl	13.126	172	9646	0.306	ng	0.00
27) Fluoranthene-d10	19.269	212	13887	0.375	ng	0.00
31) Terphenyl-d14	19.863	244	8346	0.380	ng	0.00

Target Compounds	Qvalue

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

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