

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031722\
 Data File : BN018996.D
 Acq On : 17 Mar 2022 16:06
 Operator : CG/JU
 Sample : PB143256BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143256BL

Quant Time: Mar 17 17:08:59 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 17:08:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.861	152	5234	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	13124	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	7681	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	12842	0.400	ng	0.00
29) Chrysene-d12	21.422	240	8597	0.400	ng	# 0.00
35) Perylene-d12	23.796	264	6170	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.413	112	3659	0.312	ng	0.00
5) Phenol-d6	7.016	99	3286	0.267	ng	0.00
8) Nitrobenzene-d5	9.014	82	2637	0.279	ng	0.00
11) 2-Methylnaphthalene-d10	12.253	152	5805	0.267	ng	0.00
14) 2,4,6-Tribromophenol	15.995	330	500	0.135	ng	0.01
15) 2-Fluorobiphenyl	13.127	172	8681	0.282	ng	0.00
27) Fluoranthene-d10	19.269	212	11809	0.310	ng	0.00
31) Terphenyl-d14	19.864	244	7825	0.371	ng	0.00

Target Compounds Qvalue

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

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