

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031722\
 Data File : BN019002.D
 Acq On : 17 Mar 2022 19:43
 Operator : CG/JU
 Sample : PB143257BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143257BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/18/2022
 Supervised By : Jagrut Upadhyay 03/18/2022

Quant Time: Mar 18 02:00:29 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 01:58:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.854	152	5198	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	12878	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	7567	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	12836	0.400	ng	0.00
29) Chrysene-d12	21.422	240	9074	0.400	ng	# 0.00
35) Perylene-d12	23.793	264	6481	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.406	112	3554	0.305	ng	0.00
5) Phenol-d6	7.017	99	3023	0.247	ng	0.00
8) Nitrobenzene-d5	9.014	82	2532m	0.273	ng	0.00
11) 2-Methylnaphthalene-d10	12.253	152	5735	0.268	ng	0.00
14) 2,4,6-Tribromophenol	15.984	330	468	0.128	ng	0.00
15) 2-Fluorobiphenyl	13.127	172	8502	0.280	ng	0.00
27) Fluoranthene-d10	19.270	212	12334	0.324	ng	0.00
31) Terphenyl-d14	19.864	244	7847	0.352	ng	0.00

Target Compounds Qvalue

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

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