

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032222\
 Data File : BN019109.D
 Acq On : 24 Mar 2022 03:22
 Operator : CG/JU
 Sample : PB143552BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143552BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/24/2022
 Supervised By : Jagrut Upadhyay 03/24/2022

Quant Time: Mar 24 05:33:12 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	4462	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	10286	0.400	ng	# 0.00
13) Acenaphthene-d10	14.485	164	5734	0.400	ng	-0.01
19) Phenanthrene-d10	17.236	188	9986	0.400	ng	0.00
29) Chrysene-d12	21.422	240	8572	0.400	ng	# 0.00
35) Perylene-d12	23.790	264	7580	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.406	112	3164	0.316	ng	0.00
5) Phenol-d6	7.024	99	2350m	0.224	ng	0.01
8) Nitrobenzene-d5	9.025	82	2022	0.273	ng	0.01
11) 2-Methylnaphthalene-d10	12.257	152	4893	0.287	ng	0.00
14) 2,4,6-Tribromophenol	15.995	330	377	0.136	ng	0.01
15) 2-Fluorobiphenyl	13.127	172	6830	0.297	ng	0.00
27) Fluoranthene-d10	19.265	212	11641	0.393	ng	0.00
31) Terphenyl-d14	19.860	244	6858	0.326	ng	0.00

Target Compounds Qvalue

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

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