

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092622\
 Data File : BN021912.D
 Acq On : 26 Sep 2022 11:00
 Operator : CG/JU
 Sample : PB147729BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147729BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/27/2022
 Supervised By : Jagrut Upadhyay 09/27/2022

Quant Time: Sep 26 17:36:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 14 01:55:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.047	152	4429	0.400	ng	-0.01
7) Naphthalene-d8	10.877	136	12807	0.400	ng	-0.01
13) Acenaphthene-d10	14.707	164	6017	0.400	ng	0.00
19) Phenanthrene-d10	17.467	188	10729	0.400	ng	# 0.01
29) Chrysene-d12	21.647	240	7583	0.400	ng	# 0.00
35) Perylene-d12	24.160	264	6293	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.577	112	4221	0.364	ng	0.00
5) Phenol-d6	7.209	99	4636	0.316	ng	0.00
8) Nitrobenzene-d5	9.233	82	3666m	0.346	ng	0.00
11) 2-Methylnaphthalene-d10	12.471	152	9712	0.301	ng	0.00
14) 2,4,6-Tribromophenol	16.220	330	534	0.194	ng	0.01
15) 2-Fluorobiphenyl	13.339	172	9005	0.367	ng	0.00
27) Fluoranthene-d10	19.490	212	10443	0.362	ng	0.00
31) Terphenyl-d14	20.080	244	6689	0.386	ng	0.00

Target Compounds Qvalue

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

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