

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070222\
 Data File : BN020622.D
 Acq On : 02 Jul 2022 10:51
 Operator : CG/JU
 Sample : PB145995BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB145995BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 07/05/2022
 Supervised By : Jagrut Upadhyay 07/05/2022

Quant Time: Jul 04 02:38:35 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN062622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 02 03:10:12 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.833	152	2992	0.400	ng	0.00
7) Naphthalene-d8	10.637	136	8803	0.400	ng	# 0.01
13) Acenaphthene-d10	14.474	164	5267	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	10829	0.400	ng	0.01
29) Chrysene-d12	21.440	240	8423	0.400	ng	# 0.00
35) Perylene-d12	23.790	264	7246	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.413	112	2367	0.285	ng	0.00
5) Phenol-d6	7.045	99	2191	0.221	ng	0.01
8) Nitrobenzene-d5	9.025	82	1519m	0.213	ng	0.03
11) 2-Methylnaphthalene-d10	12.235	152	4612	0.306	ng	0.00
14) 2,4,6-Tribromophenol	15.995	330	265	0.134	ng	0.02
15) 2-Fluorobiphenyl	13.116	172	5890	0.271	ng	0.01
27) Fluoranthene-d10	19.270	212	10023	0.312	ng	0.00
31) Terphenyl-d14	19.872	244	6799	0.338	ng	0.00

Target Compounds Qvalue

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

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