

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081622\
 Data File : BN021266.D
 Acq On : 16 Aug 2022 22:51
 Operator : CG/JU
 Sample : PB146962BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB146962BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/17/2022
 Supervised By : Jagrut Upadhyay 08/17/2022

Quant Time: Aug 17 05:15:51 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN081622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 16 09:19:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	2368	0.400	ng	0.01
7) Naphthalene-d8	10.603	136	7411	0.400	ng	# 0.02
13) Acenaphthene-d10	14.439	164	4733	0.400	ng	0.01
19) Phenanthrene-d10	17.203	188	8727	0.400	ng	0.02
29) Chrysene-d12	21.395	240	5927	0.400	ng	# 0.00
35) Perylene-d12	23.729	264	4849	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.386	112	2087	0.353	ng	0.00
5) Phenol-d6	7.068	99	1382m	0.200	ng	0.09
8) Nitrobenzene-d5	9.033	82	1411m	0.338	ng	0.08
11) 2-Methylnaphthalene-d10	12.210	152	4674	0.362	ng	0.03
14) 2,4,6-Tribromophenol	15.972	330	257	0.166	ng	0.03
15) 2-Fluorobiphenyl	13.082	172	5035	0.267	ng	0.02
27) Fluoranthene-d10	19.235	212	9287	0.392	ng	0.00
31) Terphenyl-d14	19.830	244	5994	0.371	ng	0.00

Target Compounds Qvalue

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

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