

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091522\
 Data File : BN021780.D
 Acq On : 15 Sep 2022 23:34
 Operator : CG/JU
 Sample : PB147657BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147657BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/16/2022
 Supervised By : Jagrut Upadhyay 09/16/2022

Quant Time: Sep 16 06:27:52 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	8.054	152	4868	0.400	ng	0.00
7) Naphthalene-d8	10.888	136	14185	0.400	ng	0.00
13) Acenaphthene-d10	14.707	164	7494	0.400	ng	0.00
19) Phenanthrene-d10	17.456	188	16124	0.400	ng	0.00
29) Chrysene-d12	21.647	240	10663	0.400	ng	# 0.00
35) Perylene-d12	24.154	264	8482	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.584	112	4887	0.384	ng	0.00
5) Phenol-d6	7.209	99	5731	0.355	ng	0.00
8) Nitrobenzene-d5	9.233	82	4068m	0.347	ng	0.00
11) 2-Methylnaphthalene-d10	12.475	152	12631	0.353	ng	0.00
14) 2,4,6-Tribromophenol	16.209	330	709	0.207	ng	0.00
15) 2-Fluorobiphenyl	13.339	172	11423	0.373	ng	0.00
27) Fluoranthene-d10	19.490	212	16192	0.373	ng	0.00
31) Terphenyl-d14	20.080	244	10355	0.425	ng	0.00

Target Compounds Qvalue

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

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