

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091422\
 Data File : BN021737.D
 Acq On : 13 Sep 2022 23:06
 Operator : CG/JU
 Sample : PB147570BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB147570BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/15/2022
 Supervised By : Jagrut Upadhyay 09/15/2022

Quant Time: Sep 14 02:12:17 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	5661	0.400	ng	0.00
7) Naphthalene-d8	10.887	136	16432	0.400	ng	0.00
13) Acenaphthene-d10	14.707	164	8718	0.400	ng	0.00
19) Phenanthrene-d10	17.456	188	19189	0.400	ng	0.00
29) Chrysene-d12	21.647	240	12406	0.400	ng	# 0.00
35) Perylene-d12	24.157	264	9663	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.584	112	5615	0.379	ng	0.00
5) Phenol-d6	7.209	99	6748	0.359	ng	0.00
8) Nitrobenzene-d5	9.232	82	4637m	0.341	ng	0.00
11) 2-Methylnaphthalene-d10	12.474	152	15059	0.363	ng	0.00
14) 2,4,6-Tribromophenol	16.208	330	959	0.241	ng	0.00
15) 2-Fluorobiphenyl	13.339	172	13796	0.388	ng	0.00
27) Fluoranthene-d10	19.489	212	19258	0.373	ng	0.00
31) Terphenyl-d14	20.080	244	10775	0.380	ng	0.00

Target Compounds Qvalue

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

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