

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120822\
 Data File : BN023101.D
 Acq On : 08 Dec 2022 19:30
 Operator : CG/JU
 Sample : PB149367BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB149367BL

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 12/13/2022
 Supervised By :Jagrut Upadhyay 12/13/2022

Quant Time: Dec 09 08:06:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN120822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 07:44:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.021	152	7246	0.400	ng	0.00
7) Naphthalene-d8	10.840	136	21512	0.400	ng	0.00
13) Acenaphthene-d10	14.666	164	11376	0.400	ng	0.00
19) Phenanthrene-d10	17.414	188	25753	0.400	ng	0.00
29) Chrysene-d12	21.589	240	18709	0.400	ng	# 0.00
35) Perylene-d12	24.056	264	16350	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.558	112	6037	0.447	ng	0.00
5) Phenol-d6	7.176	99	7809	0.455	ng	0.00
8) Nitrobenzene-d5	9.185	82	4897m	0.346	ng	0.00
11) 2-Methylnaphthalene-d10	12.427	152	11980	0.328	ng	0.00
14) 2,4,6-Tribromophenol	16.148	330	777	0.188	ng	0.00
15) 2-Fluorobiphenyl	13.298	172	16977	0.374	ng	0.00
27) Fluoranthene-d10	19.439	212	20438	0.339	ng	0.00
31) Terphenyl-d14	20.031	244	13131	0.432	ng	0.00

Target Compounds Qvalue

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

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