

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120222\
 Data File : BN023047.D
 Acq On : 02 Dec 2022 14:31
 Operator : CG/JU
 Sample : PB149383BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB149383BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 12/05/2022
 Supervised By : Jagrut Upadhyay 12/05/2022

Quant Time: Dec 05 00:33:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN112522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 02 02:30:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.049	152	6802	0.400	ng	0.00
7) Naphthalene-d8	10.861	136	17940	0.400	ng	-0.01
13) Acenaphthene-d10	14.687	164	8641	0.400	ng	0.00
19) Phenanthrene-d10	17.439	188	16742	0.400	ng	# 0.00
29) Chrysene-d12	21.625	240	9396	0.400	ng	# 0.00
35) Perylene-d12	24.106	264	6766	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.586	112	6472	0.366	ng	0.00
5) Phenol-d6	7.190	99	8082	0.359	ng	-0.01
8) Nitrobenzene-d5	9.206	82	5115m	0.367	ng	-0.01
11) 2-Methylnaphthalene-d10	12.450	152	11856	0.305	ng	0.00
14) 2,4,6-Tribromophenol	16.173	330	1249	0.299	ng	-0.01
15) 2-Fluorobiphenyl	13.319	172	15626	0.395	ng	0.00
27) Fluoranthene-d10	19.464	212	13803	0.304	ng	0.00
31) Terphenyl-d14	20.058	244	9382	0.425	ng	0.00

Target Compounds Qvalue

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

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