

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031122\
 Data File : BN018907.D
 Acq On : 10 Mar 2022 17:50
 Operator : CG/JU
 Sample : PB143229BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143229BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/11/2022
 Supervised By : Jagrut Upadhyay 03/11/2022

Quant Time: Mar 11 03:04:58 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN030322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 11 03:02:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.912	152	4446	0.400	ng	0.00
7) Naphthalene-d8	10.712	136	12348	0.400	ng	# 0.00
13) Acenaphthene-d10	14.538	164	7789	0.400	ng	0.00
19) Phenanthrene-d10	17.277	188	18907	0.400	ng	0.00
29) Chrysene-d12	21.458	240	17340	0.400	ng	0.00
35) Perylene-d12	23.846	264	15162	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.464	112	3104	0.319	ng	0.00
5) Phenol-d6	7.067	99	3449	0.296	ng	0.00
8) Nitrobenzene-d5	9.067	82	2440m	0.271	ng	0.01
11) 2-Methylnaphthalene-d10	12.302	152	5488	0.273	ng	0.00
14) 2,4,6-Tribromophenol	16.025	330	651	0.181	ng	0.00
15) 2-Fluorobiphenyl	13.169	172	9055	0.271	ng	0.00
27) Fluoranthene-d10	19.303	212	15182	0.273	ng	0.00
31) Terphenyl-d14	19.898	244	10807	0.292	ng	0.00

Target Compounds Qvalue

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

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