

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042222\
 Data File : BN019521.D
 Acq On : 22 Apr 2022 14:24
 Operator : CG/JU
 Sample : PB144158BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB144158BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/25/2022
 Supervised By : Jagrut Upadhyay 04/25/2022

Quant Time: Apr 23 06:14:39 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN042122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 20 17:50:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.819	152	3864	0.400	ng	0.00
7) Naphthalene-d8	10.645	136	9729	0.400	ng	# 0.03
13) Acenaphthene-d10	14.461	164	6203	0.400	ng	0.02
19) Phenanthrene-d10	17.213	188	12159	0.400	ng	0.02
29) Chrysene-d12	21.395	240	8467	0.400	ng	# 0.01
35) Perylene-d12	23.740	264	7052	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.371	112	2983	0.360	ng	0.00
5) Phenol-d6	7.040	99	1960m	0.208	ng	0.06
8) Nitrobenzene-d5	9.097	82	1634m	0.232	ng	0.12
11) 2-Methylnaphthalene-d10	12.244	152	5252	0.333	ng	0.03
14) 2,4,6-Tribromophenol	15.982	330	483m	0.174	ng	0.04
15) 2-Fluorobiphenyl	13.114	172	6479m	0.262	ng	0.03
27) Fluoranthene-d10	19.235	212	13406	0.377	ng	0.01
31) Terphenyl-d14	19.834	244	7373	0.386	ng	0.00

Target Compounds Qvalue

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

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