

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120921\
 Data File : BN017853.D
 Acq On : 10 Dec 2021 03:31
 Operator : CG/JU
 Sample : PB141233BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB141233BL

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021
 Supervised By :Yogesh Patel 12/15/2021

Quant Time: Dec 10 06:51:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN120721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 08 01:18:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.715	152	20337	0.400	ng	0.00
7) Naphthalene-d8	10.507	136	72062	0.400	ng	0.01
13) Acenaphthene-d10	14.347	164	45938	0.400	ng	0.00
19) Phenanthrene-d10	17.117	188	82076	0.400	ng	# 0.02
29) Chrysene-d12	21.303	240	66207	0.400	ng	0.02
35) Perylene-d12	23.596	264	45828	0.400	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.345	112	12311	0.251	ng	0.02
5) Phenol-d6	6.912	99	8395	0.181	ng	0.00
8) Nitrobenzene-d5	8.859	82	14036	0.293	ng	0.00
11) 2-Methylnaphthalene-d10	12.103	152	36488	0.254	ng	0.01
14) 2,4,6-Tribromophenol	15.882	330	1235m	0.051	ng	0.04
15) 2-Fluorobiphenyl	12.990	172	52728m	0.313	ng	0.02
27) Fluoranthene-d10	19.139	212	91412	0.333	ng	0.00
31) Terphenyl-d14	19.745	244	71704	0.504	ng	0.00

Target Compounds Qvalue

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

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