

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012522\
 Data File : BN018465.D
 Acq On : 25 Jan 2022 18:37
 Operator : CG/JU
 Sample : PB142256BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB142256BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/26/2022
 Supervised By : Jagrut Upadhyay 01/26/2022

Quant Time: Jan 26 04:07:43 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 18:25:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.956	152	34246	0.400	ng	0.00
7) Naphthalene-d8	10.761	136	150745	0.400	ng	0.00
13) Acenaphthene-d10	14.573	164	78521	0.400	ng	0.00
19) Phenanthrene-d10	17.309	188	144834	0.400	ng	# 0.00
29) Chrysene-d12	21.482	240	96453	0.400	ng	-0.01
35) Perylene-d12	23.888	264	68017	0.400	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.531	112	32085	0.378	ng	0.02
5) Phenol-d6	7.116	99	34141	0.370	ng	0.00
8) Nitrobenzene-d5	9.101	82	33097m	0.381	ng	-0.01
11) 2-Methylnaphthalene-d10	12.340	152	89452	0.354	ng	0.00
14) 2,4,6-Tribromophenol	16.056	330	4883	0.206	ng	0.00
15) 2-Fluorobiphenyl	13.203	172	125147	0.444	ng	-0.01
27) Fluoranthene-d10	19.331	212	149812	0.343	ng	-0.01
31) Terphenyl-d14	19.926	244	134835	0.450	ng	-0.01

Target Compounds Qvalue

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

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