

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070221\
 Data File : BN015313.D
 Acq On : 03 Jul 2021 01:21
 Operator : CG/JU
 Sample : M2803-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW202D1-20210619

Manual Integrations
 APPROVED

mohammad
 7/6/2021 10:03:38 AM

Quant Time: Jul 03 02:15:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN063021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 01 09:06:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.735	152	22704	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	104296	0.40	ng	0.00
13) Acenaphthene-d10	14.345	164	56598	0.40	ng	0.00
19) Phenanthrene-d10	17.091	188	113016	0.40	ng	# 0.00
29) Chrysene-d12	21.291	240	83759	0.40	ng	#-0.01
35) Perylene-d12	23.532	264	68478	0.40	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.365	112	9902	0.17	ng	0.00
5) Phenol-d6	6.905	99	6703	0.09	ng	0.00
8) Nitrobenzene-d5	8.860	82	28561	0.49	ng	0.00
11) 2-Methylnaphthalene-d10	12.083	152	60294	0.35	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	7390	0.52	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	85817	0.37	ng	0.00
27) Fluoranthene-d10	19.127	212	130766	0.39	ng	0.00
31) Terphenyl-d14	19.728	244	105326	0.54	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.355	88	4101m	0.39	ng	
24) Pentachlorophenol	16.750	266	80	0.25	ng	# 89
34) Bis(2-ethylhexyl)phtha...	21.206	149	28489	0.43	ng	# 87

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

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