

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN063021\
 Data File : BN015261.D
 Acq On : 01 Jul 2021 15:42
 Operator : CG/JU
 Sample : M2803-11
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW202D-20210619

Manual Integrations
 APPROVED

mohammad
 7/2/2021 7:57:03 AM

Quant Time: Jul 01 16:17:55 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	7.734	152	24226	0.400	ng	# 0.00
7) Naphthalene-d8	10.495	136	110916	0.400	ng	0.00
13) Acenaphthene-d10	14.345	164	57768	0.400	ng	0.00
19) Phenanthrene-d10	17.090	188	113505	0.400	ng	0.00
29) Chrysene-d12	21.291	240	92724	0.400	ng	#-0.01
35) Perylene-d12	23.539	264	75466	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.365	112	9114	0.148	ng	0.00
5) Phenol-d6	6.914	99	6050	0.078	ng	0.01
8) Nitrobenzene-d5	8.873	82	27611	0.444	ng	0.00
11) 2-Methylnaphthalene-d10	12.087	152	62816	0.342	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	6028	0.462	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	89210	0.380	ng	0.00
27) Fluoranthene-d10	19.127	212	129690	0.387	ng	0.00
31) Terphenyl-d14	19.733	244	104504	0.482	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.364	88	7102m	0.625	ng	Qvalue
24) Pentachlorophenol	16.750	266	239	0.260	ng	96
34) Bis(2-ethylhexyl)phtha...	21.217	149	22724	0.350	ng	# 86

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

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