

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070621\
 Data File : BN015342.D
 Acq On : 06 Jul 2021 17:24
 Operator : CG/JU
 Sample : M2803-22DL 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D1-20210621DL

Manual Integrations
 APPROVED

mohammad
 7/7/2021 8:35:59 AM

Quant Time: Jul 06 17:52:37 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	18395	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	87836	0.40	ng	# 0.00
13) Acenaphthene-d10	14.344	164	52532	0.40	ng	0.00
19) Phenanthrene-d10	17.090	188	107843	0.40	ng	0.00
29) Chrysene-d12	21.301	240	90528	0.40	ng	# 0.00
35) Perylene-d12	23.556	264	52529	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.374	112	3785	0.08	ng	0.01
5) Phenol-d6	6.914	99	2687	0.05	ng	0.01
8) Nitrobenzene-d5	8.873	82	11567	0.24	ng	0.00
11) 2-Methylnaphthalene-d10	12.087	152	25133	0.17	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	4656	0.43	ng	0.00
15) 2-Fluorobiphenyl	12.974	172	36575	0.17	ng	0.01
27) Fluoranthene-d10	19.132	212	60384	0.19	ng	0.00
31) Terphenyl-d14	19.738	244	50191	0.24	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.345	88	9783	1.13	ng	# 62
6) bis(2-Chloroethyl)ether	7.172	93	1847m	0.03	ng	

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

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