

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070221\
 Data File : BN015311.D
 Acq On : 03 Jul 2021 00:10
 Operator : CG/JU
 Sample : M2803-22
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D1-20210621

Manual Integrations
 APPROVED

mohammad
 7/6/2021 10:03:30 AM

Quant Time: Jul 03 00:54:01 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	22655	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	102759	0.40	ng	0.00
13) Acenaphthene-d10	14.345	164	56655	0.40	ng	0.00
19) Phenanthrene-d10	17.091	188	110040	0.40	ng	# 0.00
29) Chrysene-d12	21.291	240	80002	0.40	ng	-0.01
35) Perylene-d12	23.529	264	62201	0.40	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.365	112	9629	0.17	ng	0.00
5) Phenol-d6	6.905	99	6561	0.09	ng	0.00
8) Nitrobenzene-d5	8.861	82	28230	0.49	ng	0.00
11) 2-Methylnaphthalene-d10	12.083	152	62904	0.37	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	7002	0.51	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	89908	0.39	ng	0.00
27) Fluoranthene-d10	19.127	212	134011	0.41	ng	0.00
31) Terphenyl-d14	19.728	244	103277	0.55	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.355	88	55037m	5.18	ng	
6) bis(2-Chloroethyl)ether	7.163	93	5289	0.07	ng	# 85
24) Pentachlorophenol	16.750	266	59	0.25	ng	87
34) Bis(2-ethylhexyl)phtha...	21.206	149	22580	0.38	ng	# 87

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

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