

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070721\
 Data File : BN015369.D
 Acq On : 07 Jul 2021 12:16
 Operator : CG/JU
 Sample : M2803-22 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D1-20210621

Manual Integrations
 APPROVED

mohammad
 7/8/2021 9:13:58 AM

Quant Time: Jul 07 12:48:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN063021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 07 11:11:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.725	152	22191	0.40	ng	# 0.00
7) Naphthalene-d8	10.483	136	101133	0.40	ng	# 0.00
13) Acenaphthene-d10	14.332	164	55471	0.40	ng	0.00
19) Phenanthrene-d10	17.078	188	110077	0.40	ng	0.00
29) Chrysene-d12	21.291	240	98486	0.40	ng	# 0.00
35) Perylene-d12	23.536	264	91417	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.356	112	4187	0.07	ng	0.02
5) Phenol-d6	6.905	99	2806	0.04	ng	0.00
8) Nitrobenzene-d5	8.860	82	12799	0.23	ng	0.00
11) 2-Methylnaphthalene-d10	12.078	152	27474	0.16	ng	0.00
14) 2,4,6-Tribromophenol	15.829	330	3300	0.36	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	39497	0.18	ng	0.00
27) Fluoranthene-d10	19.122	212	60842	0.19	ng	0.00
31) Terphenyl-d14	19.728	244	48609	0.21	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.345	88	21558m	2.07	ng	Qvalue
6) bis(2-Chloroethyl)ether	7.163	93	2282	0.03	ng	# 86
34) Bis(2-ethylhexyl)phtha...	21.217	149	10129	0.23	ng	# 88

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

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