

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070221\
 Data File : BN015309.D
 Acq On : 02 Jul 2021 22:58
 Operator : CG/JU
 Sample : M2803-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT180D2-20210618

Manual Integrations
 APPROVED

mohammad
 7/6/2021 10:03:25 AM

Quant Time: Jul 03 00:53:42 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	21228	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	96616	0.40	ng	0.00
13) Acenaphthene-d10	14.345	164	51030	0.40	ng	0.00
19) Phenanthrene-d10	17.091	188	97065	0.40	ng	0.00
29) Chrysene-d12	21.291	240	67636	0.40	ng	#-0.01
35) Perylene-d12	23.536	264	52766	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.365	112	8904	0.16	ng	0.00
5) Phenol-d6	6.905	99	5892	0.09	ng	0.00
8) Nitrobenzene-d5	8.860	82	25600	0.47	ng	0.00
11) 2-Methylnaphthalene-d10	12.083	152	55105	0.34	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	5036	0.45	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	81209	0.39	ng	0.00
27) Fluoranthene-d10	19.127	212	104749	0.37	ng	0.00
31) Terphenyl-d14	19.728	244	89439	0.57	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.355	88	44731m	4.49	ng	
6) bis(2-Chloroethyl)ether	7.163	93	3140	0.05	ng	# 84
24) Pentachlorophenol	16.750	266	111	0.25	ng	88
34) Bis(2-ethylhexyl)phtha...	21.206	149	23765	0.43	ng	# 87

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

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