

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070221\
 Data File : BN015322.D
 Acq On : 03 Jul 2021 07:53
 Operator : CG/JU
 Sample : M2838-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE103D3-20210621

Manual Integrations
 APPROVED

mohammad
 7/6/2021 10:03:50 AM

Quant Time: Jul 05 05:48:31 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	25344	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	116523	0.40	ng	0.00
13) Acenaphthene-d10	14.345	164	63972	0.40	ng	0.00
19) Phenanthrene-d10	17.091	188	123966	0.40	ng	# 0.00
29) Chrysene-d12	21.291	240	83488	0.40	ng	-0.01
35) Perylene-d12	23.532	264	63162	0.40	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.365	112	10035	0.16	ng	0.00
5) Phenol-d6	6.905	99	6791	0.08	ng	0.00
8) Nitrobenzene-d5	8.861	82	28548	0.44	ng	0.00
11) 2-Methylnaphthalene-d10	12.083	152	64782	0.34	ng	0.00
14) 2,4,6-Tribromophenol	15.829	330	6991	0.47	ng	-0.01
15) 2-Fluorobiphenyl	12.962	172	95799	0.37	ng	0.00
27) Fluoranthene-d10	19.123	212	124298	0.34	ng	-0.01
31) Terphenyl-d14	19.728	244	98770	0.51	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.355	88	7484m	0.63	ng	
24) Pentachlorophenol	16.750	266	109	0.25	ng	# 77
34) Bis(2-ethylhexyl)phtha...	21.206	149	31880	0.46	ng	# 86

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

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