

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN062220\
 Data File : VN062006.D
 Acq On : 22 Jun 2020 17:55
 Operator : JC/MD
 Sample : L3033-03DL 4X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SS043MW011-200615DL

Manual Integrations
 APPROVED

MMDadoda
 6/23/2020 12:47:41 PM

Quant Time: Jun 23 06:04:57 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 17 01:19:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	147397	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	274018	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	275070	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	104918	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	104103	52.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.32%	
35) Dibromofluoromethane	7.55	113	88688	50.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.18%	
50) Toluene-d8	10.06	98	348567	50.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.16%	
62) 4-Bromofluorobenzene	12.38	95	117790	49.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.66%	
Target Compounds						
6) Chloroethane	2.67	64	2086	1.264	ug/l	92
8) Diethyl Ether	3.38	74	1845	1.884	ug/l	73
29) Tetrahydrofuran	7.17	42	1565m	2.791	ug/l	
40) Benzene	8.00	78	5827	0.770	ug/l	93
49) 1,4-Dioxane	9.17	88	6710	195.390	ug/l	95
52) Toluene	10.12	92	4592	0.966	ug/l	92
65) Chlorobenzene	11.41	112	360149	62.050	ug/l	100
87) 1,3-Dichlorobenzene	13.26	146	1596	0.465	ug/l	88
88) 1,4-Dichlorobenzene	13.34	146	12629	3.596	ug/l	88
91) 1,2-Dichlorobenzene	13.63	146	4545	1.376	ug/l	95

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

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