

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061120\
 Data File : VN061757.D
 Acq On : 11 Jun 2020 19:10
 Operator : JC/MD
 Sample : L2988-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RE132D1-20200608

Manual Integrations
 APPROVED

MMDadoda
 6/12/2020 5:46:17 PM

Quant Time: Jun 12 06:23:15 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N052820W.M
 Quant Title : SW846 8260
 QLast Update : Fri May 29 02:14:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	279689	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	475365	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	465888	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	203026	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	159408	47.84	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.68%	
35) Dibromofluoromethane	7.55	113	138674	57.51	ug/l	0.00
Spiked Amount	50.000		Recovery	=	115.02%	
50) Toluene-d8	10.06	98	574992	51.45	ug/l	0.00
Spiked Amount	50.000		Recovery	=	102.90%	
62) 4-Bromofluorobenzene	12.38	95	193757	47.52	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.04%	
Target Compounds						
2) Dichlorodifluoromethane	1.83	85	1074	1.473	ug/l	81
9) 1,1,2-Trichlorotrifluoroet	3.72	101	15968	5.969	ug/l	96
12) 1,1-Dichloroethene	3.71	96	2829	0.990	ug/l #	66
19) Methyl tert-butyl Ether	5.00	73	8617	0.885	ug/l	98
22) Diisopropyl ether	5.90	45	4121m	0.500	ug/l	
27) cis-1,2-Dichloroethene	6.78	96	6390	1.751	ug/l	94
30) Chloroform	7.33	83	10401	1.800	ug/l	100
44) Trichloroethene	8.80	130	400925	100.473	ug/l	97
64) Tetrachloroethene	10.61	164	11578	2.640	ug/l	95

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

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