

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061820\
 Data File : VN061899.D
 Acq On : 18 Jun 2020 14:50
 Operator : JC/MD
 Sample : L3011-17
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RE104D2-20200612

Quant Time: Jun 19 13:08:44 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	138441	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	251617	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	239808	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	89450	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	90867	48.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.88%	
35) Dibromofluoromethane	7.55	113	81137	50.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.80%	
50) Toluene-d8	10.06	98	311673	49.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.50%	
62) 4-Bromofluorobenzene	12.38	95	101207	46.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.24%	
Target Compounds						
12) 1,1-Dichloroethene	3.70	96	1997	1.420	ug/l	80
16) Acetone	3.79	43	2246	4.095	ug/l	99
27) cis-1,2-Dichloroethene	6.78	96	90049	48.303	ug/l	89
30) Chloroform	7.33	83	14654	4.906	ug/l	92
42) 1,2-Dichloroethane	8.08	62	4787	2.019	ug/l	84
44) Trichloroethene	8.80	130	568447	310.689	ug/l	98

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

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