

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN013120\
 Data File : VN059949.D
 Acq On : 31 Jan 2020 15:42
 Operator : JC/MD
 Sample : L1333-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1-2-COMP

Quant Time: Feb 03 03:22:40 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011320W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jan 18 03:09:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	195130	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	300054	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	269743	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	119817	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	107546	49.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.26%	
35) Dibromofluoromethane	7.57	113	90574	50.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.24%	
50) Toluene-d8	10.08	98	358757	51.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.50%	
62) 4-Bromofluorobenzene	12.40	95	122052	48.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.66%	
Target Compounds						
16) Acetone	3.79	43	13413	8.477	ug/l	99
25) 2-Butanone	6.81	43	4872	3.970	ug/l	98
43) Isopropyl Acetate	8.14	43	48274	11.452	ug/l #	90
51) 4-Methyl-2-Pentanone	9.97	43	2871	1.178	ug/l	97
68) m/p-Xylenes	11.61	106	14246	3.918	ug/l	96
69) o-Xylene	11.94	106	5424	1.524	ug/l	98

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

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