

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091319\
 Data File : VN058000.D
 Acq On : 13 Sep 2019 3:05
 Operator : JC/SP
 Sample : K4644-30
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T11-E2-(4)

Quant Time: Sep 13 06:58:00 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	267032	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	460664	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	421710	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	215855	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	317353	58.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	117.36%	
35) Dibromofluoromethane	7.58	113	214956	50.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.12%	
50) Toluene-d8	10.09	98	830693	50.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.64%	
62) 4-Bromofluorobenzene	12.41	95	275947	43.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.72%	
Target Compounds						
16) Acetone	3.80	43	15986	9.016	ug/l	100
18) Methyl Acetate	4.30	43	74556	15.771	ug/l	99
20) Methylene Chloride	4.53	84	7803	1.108	ug/l #	84
43) Isopropyl Acetate	8.16	43	20596	2.219	ug/l #	78

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

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