

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091419\
 Data File : VN058030.D
 Acq On : 13 Sep 2019 20:29
 Operator : JC/SP
 Sample : K4680-37
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T3-4(0)

Quant Time: Sep 14 05:37:43 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	288880	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	495539	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	446344	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	223566	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	328239	56.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.20%	
35) Dibromofluoromethane	7.58	113	226758	49.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.16%	
50) Toluene-d8	10.09	98	780296	43.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.88%	
62) 4-Bromofluorobenzene	12.40	95	269888	39.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	79.76%	
Target Compounds						
16) Acetone	3.80	43	15942	8.311	ug/l	91
18) Methyl Acetate	4.31	43	7108	1.390	ug/l #	71
20) Methylene Chloride	4.53	84	13624	2.563	ug/l	97
43) Isopropyl Acetate	8.16	43	81720	8.184	ug/l #	86

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

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