

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN103119\
 Data File : VN059037.D
 Acq On : 31 Oct 2019 14:17
 Operator : JC/SP
 Sample : K5615-06
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 3465

Quant Time: Nov 01 02:33:33 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N101819W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 19 00:16:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	384648	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	604087	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	552041	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	197547	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	265771	47.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.22%	
35) Dibromofluoromethane	7.57	113	197220	49.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.32%	
50) Toluene-d8	10.09	98	769195	49.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.78%	
62) 4-Bromofluorobenzene	12.41	95	242536	42.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	85.68%	
Target Compounds						
8) Diethyl Ether	3.39	74	4579	1.816	ug/l	97
16) Acetone	3.80	43	28182	10.109	ug/l	92
18) Methyl Acetate	4.31	43	11385	1.796	ug/l	92
20) Methylene Chloride	4.53	84	267868	56.274	ug/l	96
30) Chloroform	7.36	83	8854	1.014	ug/l	89
43) Isopropyl Acetate	8.15	43	87571	8.103	ug/l #	91

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

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