

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN010919\
 Data File : VN053323.D
 Acq On : 9 Jan 2019 16:38
 Operator : MD\SY
 Sample : K1100-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 8 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 WATER-DRUMS-A

Quant Time: Jan 10 03:34:11 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N010519W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jan 07 09:34:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	1111979	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1704114	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1555032	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	656962	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	538739	46.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.14%	
35) Dibromofluoromethane	7.59	113	492061	50.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.44%	
50) Toluene-d8	10.09	98	2094922	50.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.12%	
62) 4-Bromofluorobenzene	12.40	95	714619	48.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.56%	
Target Compounds						
16) Acetone	3.82	43	42129	13.012	ug/l	98
18) Methyl Acetate	4.33	43	7114	0.679	ug/l #	75
20) Methylene Chloride	4.56	84	5282	0.425	ug/l	89
25) 2-Butanone	6.84	43	9069	1.911	ug/l	100
30) Chloroform	7.37	83	195994	9.584	ug/l	100
47) Bromodichloromethane	9.40	83	9649	0.623	ug/l	94
52) Toluene	10.16	92	25702	0.862	ug/l	98

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

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