

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061019\
 Data File : VN056187.D
 Acq On : 10 Jun 2019 23:44
 Operator : JC/SP
 Sample : K3267-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP19-CHY05

Quant Time: Jun 11 04:46:52 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 04 11:02:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	310715	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	473809	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	466852	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	159399	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	169665	54.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.98%	
35) Dibromofluoromethane	7.59	113	155584	53.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.04%	
50) Toluene-d8	10.09	98	608190	54.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.74%	
62) 4-Bromofluorobenzene	12.40	95	201452	52.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.92%	
Target Compounds						
16) Acetone	3.82	43	10870	7.367	ug/l	99
18) Methyl Acetate	4.34	43	6377	1.641	ug/l #	71
27) cis-1,2-Dichloroethene	6.82	96	9330	2.840	ug/l	93
44) Trichloroethene	8.83	130	13314	3.744	ug/l	94

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

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