

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN041919\
 Data File : VN055151.D
 Acq On : 19 Apr 2019 20:47
 Operator : JC/SP
 Sample : K2201-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PL-01-041819-C

Quant Time: Apr 20 04:58:10 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N041519W.M
 Quant Title : SW846 8260
 QLast Update : Mon Apr 15 12:28:28 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	542249	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	861406	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	676619	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	195059	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	351702	48.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.58%	
35) Dibromofluoromethane	7.59	113	277500	48.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.40%	
50) Toluene-d8	10.09	98	1008699	48.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.92%	
62) 4-Bromofluorobenzene	12.40	95	270584	40.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	80.80%	
Target Compounds						
16) Acetone	3.83	43	14840	7.586	ug/l #	80
20) Methylene Chloride	4.55	84	513212	74.625	ug/l	98
30) Chloroform	7.37	83	12871	1.149	ug/l	97
43) Isopropyl Acetate	8.17	43	13086	1.285	ug/l	97

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

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