

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031320\
 Data File : VN060511.D
 Acq On : 13 Mar 2020 18:28
 Operator : JC/MD
 Sample : L1907-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-9-SA

Quant Time: Mar 16 09:20:43 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N022720W.M
 Quant Title : SW846 8260
 QLast Update : Thu Feb 27 13:52:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	201041	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	333177	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	316626	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	149715	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	131470	50.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.94%	
35) Dibromofluoromethane	7.57	113	99370	64.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	128.50%	
50) Toluene-d8	10.08	98	409501	52.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.64%	
62) 4-Bromofluorobenzene	12.40	95	156609	52.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.56%	
Target Compounds						
16) Acetone	3.79	43	7636	9.167	ug/l	96
20) Methylene Chloride	4.53	84	7587	3.317	ug/l	97
43) Isopropyl Acetate	8.14	43	64118	12.181	ug/l #	90
51) 4-Methyl-2-Pentanone	9.97	43	4995	1.674	ug/l	94

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

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