

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN073120\
 Data File : VN062719.D
 Acq On : 1 Aug 2020 5:36
 Operator : JC/MD
 Sample : L3518-33
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-10

Quant Time: Aug 01 06:57:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 21 18:48:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	130316	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	242280	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	244164	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	91282	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	117615	58.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	116.20%	
35) Dibromofluoromethane	7.55	113	80486	50.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.14%	
50) Toluene-d8	10.06	98	317866	50.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.12%	
62) 4-Bromofluorobenzene	12.38	95	118794	51.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.34%	
Target Compounds						
16) Acetone	3.78	43	12943	16.990	ug/l	95
18) Methyl Acetate	4.28	43	5119	2.654	ug/l #	81
20) Methylene Chloride	4.51	84	7498	4.104	ug/l	92
43) Isopropyl Acetate	8.13	43	44069	9.820	ug/l #	89

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

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