

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN072720\
 Data File : VN062619.D
 Acq On : 27 Jul 2020 23:26
 Operator : JC/MD
 Sample : L3439-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 20200440-3001C

Quant Time: Jul 28 07:58:02 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	135145	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	242254	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	242755	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	107740	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	116329	55.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.82%	
35) Dibromofluoromethane	7.55	113	81887	51.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.92%	
50) Toluene-d8	10.06	98	316100	49.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.58%	
62) 4-Bromofluorobenzene	12.38	95	121759	52.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.90%	
Target Compounds						
16) Acetone	3.78	43	80436	101.815	ug/l	100
20) Methylene Chloride	4.50	84	10495	5.539	ug/l	89
25) 2-Butanone	6.79	43	19288	16.031	ug/l	97
43) Isopropyl Acetate	8.13	43	15496	3.453	ug/l #	89

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

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