

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN063020\
 Data File : VN062270.D
 Acq On : 1 Jul 2020 4:43
 Operator : JC/MD
 Sample : L3153-11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP2-A

Quant Time: Jul 01 06:11:25 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 24 15:33:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	208933	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	425100	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	416299	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	155952	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	162266	55.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.68%	
35) Dibromofluoromethane	7.55	113	130793	47.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.92%	
50) Toluene-d8	10.06	98	544083	49.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.82%	
62) 4-Bromofluorobenzene	12.38	95	180988	49.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.28%	
Target Compounds						
16) Acetone	3.78	43	78989	83.123	ug/l	98
18) Methyl Acetate	4.29	43	6622	2.841	ug/l #	79
20) Methylene Chloride	4.50	84	23697	8.625	ug/l	98
43) Isopropyl Acetate	8.12	43	35174	6.098	ug/l	98
79) 2-Chlorotoluene	12.65	91	11004	1.318	ug/l	100
82) 4-Chlorotoluene	12.75	91	11731	1.359	ug/l	97

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

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