

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN052120\
 Data File : VN061536.D
 Acq On : 21 May 2020 17:24
 Operator : JC/MD
 Sample : L2727-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CHASSIS-WASH

Quant Time: May 22 08:21:05 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
 Quant Title : SW846 8260
 QLast Update : Mon May 18 13:09:55 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	286799	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	523633	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	527567	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	249728	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	185077	43.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.48%	
35) Dibromofluoromethane	7.55	113	151031	55.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.12%	
50) Toluene-d8	10.06	98	656946	51.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.26%	
62) 4-Bromofluorobenzene	12.37	95	254007	53.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.86%	
Target Compounds						
7) Trichlorofluoromethane	2.98	101	9651	1.428	ug/l	99
16) Acetone	3.77	43	70529	55.507	ug/l	97
20) Methylene Chloride	4.50	84	18021	5.123	ug/l	90
25) 2-Butanone	6.79	43	17752	9.109	ug/l	94
30) Chloroform	7.33	83	18205	2.780	ug/l	97
37) Ethyl Acetate	6.89	43	8617	1.941	ug/l	98
43) Isopropyl Acetate	8.13	43	63558	8.207	ug/l #	91
51) 4-Methyl-2-Pentanone	9.95	43	92217	21.858	ug/l	98

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

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