

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN033120\
 Data File : VN060788.D
 Acq On : 31 Mar 2020 16:05
 Operator : JC/MD
 Sample : L2070-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BUSH-WASH

Quant Time: Apr 01 09:13:37 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 18 08:49:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	200309	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.56	114	336699	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	314392	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	145335	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	129938	47.25	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.50%	
35) Dibromofluoromethane	7.56	113	99238	48.74	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.48%	
50) Toluene-d8	10.08	98	416388	50.07	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.14%	
62) 4-Bromofluorobenzene	12.40	95	155177	50.92	ug/l	0.00
Spiked Amount	50.000		Recovery	=	101.84%	
Target Compounds						
16) Acetone	3.78	43	63683	65.381	ug/l	98
18) Methyl Acetate	4.29	43	12278	4.478	ug/l	94
20) Methylene Chloride	4.52	84	9019	3.764	ug/l	97
25) 2-Butanone	6.80	43	22349	14.179	ug/l	96
43) Isopropyl Acetate	8.14	43	13072	2.250	ug/l #	90
51) 4-Methyl-2-Pentanone	9.97	43	124321	37.604	ug/l	99

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

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