

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN041420\
 Data File : VN060941.D
 Acq On : 14 Apr 2020 18:08
 Operator : JC/MD
 Sample : L2243-13
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-31-30-COMP

Quant Time: Apr 15 09:39:47 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	145537	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.56	114	249141	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	237415	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	106817	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	97901	49.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.00%	
35) Dibromofluoromethane	7.56	113	73469	48.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.54%	
50) Toluene-d8	10.08	98	310527	50.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.92%	
62) 4-Bromofluorobenzene	12.40	95	117951	52.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.60%	
Target Compounds						
16) Acetone	3.78	43	11942	16.875	ug/l	98
20) Methylene Chloride	4.52	84	2866	1.646	ug/l	94
25) 2-Butanone	6.81	43	3427	2.993	ug/l	92
43) Isopropyl Acetate	8.14	43	48193	11.210	ug/l #	90

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

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