

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN040220\
 Data File : VN060835.D
 Acq On : 2 Apr 2020 20:28
 Operator : JC/MD
 Sample : L2104-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OIL-WATER-SEPARATOR

Quant Time: Apr 03 09:30:47 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	179133	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.56	114	297856	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	278103	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	128129	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	117680	47.86	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.72%	
35) Dibromofluoromethane	7.56	113	89484	49.68	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.36%	
50) Toluene-d8	10.08	98	370849	50.41	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.82%	
62) 4-Bromofluorobenzene	12.40	95	135967	50.43	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.86%	
Target Compounds						
16) Acetone	3.78	43	34010	39.044	ug/l	100
20) Methylene Chloride	4.52	84	6919	3.229	ug/l	99
25) 2-Butanone	6.80	43	15659	11.109	ug/l	96
43) Isopropyl Acetate	8.14	43	7892	1.536	ug/l #	89
51) 4-Methyl-2-Pentanone	9.97	43	75144	25.693	ug/l	99
52) Toluene	10.14	92	89925	17.102	ug/l	99

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

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