

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN041620\
 Data File : VN061016.D
 Acq On : 16 Apr 2020 17:04
 Operator : JC/MD
 Sample : L2259-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OIL-WATER-SEPARATOR-2

Quant Time: Apr 17 04:40:41 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N041620W.M
 Quant Title : SW846 8260
 QLast Update : Fri Apr 17 03:49:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	159564	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.56	114	271005	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	253325	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	112658	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	108553	53.17	ug/l	0.00
Spiked Amount	50.000		Recovery	=	106.34%	
35) Dibromofluoromethane	7.56	113	81972	51.79	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.58%	
50) Toluene-d8	10.08	98	337833	56.49	ug/l	0.00
Spiked Amount	50.000		Recovery	=	112.98%	
62) 4-Bromofluorobenzene	12.39	95	124415	53.08	ug/l	0.00
Spiked Amount	50.000		Recovery	=	106.16%	
Target Compounds						
16) Acetone	3.78	43	50112	55.768	ug/l	99
20) Methylene Chloride	4.52	84	17521	8.931	ug/l	93
25) 2-Butanone	6.80	43	19340	14.046	ug/l	96
30) Chloroform	7.35	83	87345	25.269	ug/l	98
43) Isopropyl Acetate	8.14	43	42323	8.570	ug/l #	91
51) 4-Methyl-2-Pentanone	9.97	43	131194	46.082	ug/l	99

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

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