

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN040219\
 Data File : VN054897.D
 Acq On : 2 Apr 2019 13:12
 Operator : JC/SP
 Sample : K2170-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 9 Sample Multiplier: 1

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

Quant Time: Apr 03 05:47:32 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N032919W.M
 Quant Title : SW846 8260
 QLast Update : Sat Mar 30 00:48:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	483214	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	814972	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	716081	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	301776	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	325465	49.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.94%	
35) Dibromofluoromethane	7.59	113	260290	47.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.82%	
50) Toluene-d8	10.09	98	998842	49.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.90%	
62) 4-Bromofluorobenzene	12.40	95	327484	48.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.50%	
Target Compounds						
					Qvalue	
16) Acetone	3.83	43	56210	36.187	ug/l #	87
20) Methylene Chloride	4.55	84	42229	7.707	ug/l #	85
25) 2-Butanone	6.86	43	24132	10.510	ug/l #	82
43) Isopropyl Acetate	8.17	43	11491	1.270	ug/l #	92
51) 4-Methyl-2-Pentanone	9.99	43	90237	17.738	ug/l #	87

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

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