

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN030320\
 Data File : VN060330.D
 Acq On : 3 Mar 2020 15:27
 Operator : JC/MD
 Sample : L1757-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EO-02-030220

Quant Time: Mar 04 02:31:13 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N022720W.M
 Quant Title : SW846 8260
 QLast Update : Thu Feb 27 13:52:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	192494	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	311536	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	296534	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	143522	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	121735	49.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.58%	
35) Dibromofluoromethane	7.57	113	88708	61.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	122.70%	
50) Toluene-d8	10.08	98	385257	52.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.28%	
62) 4-Bromofluorobenzene	12.40	95	145529	52.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.92%	
Target Compounds						
16) Acetone	3.79	43	27547	34.540	ug/l	96
20) Methylene Chloride	4.52	84	5981	2.731	ug/l #	92
43) Isopropyl Acetate	8.14	43	34420	6.993	ug/l #	91
67) Ethyl Benzene	11.51	91	12683	1.163	ug/l	98
68) m/p-Xylenes	11.61	106	17196	4.180	ug/l	97
69) o-Xylene	11.94	106	16146	4.116	ug/l	99
80) 1,3,5-Trimethylbenzene	12.73	105	26011	2.904	ug/l	99
84) 1,2,4-Trimethylbenzene	13.04	105	87734	9.705	ug/l	100

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

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