

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061820\
 Data File : VN061941.D
 Acq On : 19 Jun 2020 8:54
 Operator : JC/MD
 Sample : L2817-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SU-04-061720

Quant Time: Jun 19 12:04:23 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 17 01:19:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	143052	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	269130	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	270971	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	101941	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	97939	51.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.10%	
35) Dibromofluoromethane	7.55	113	83959	48.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.52%	
50) Toluene-d8	10.06	98	344830	50.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.88%	
62) 4-Bromofluorobenzene	12.38	95	114668	49.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.78%	
Target Compounds						
16) Acetone	3.78	43	17121	30.212	ug/l	99
20) Methylene Chloride	4.51	84	6160	3.336	ug/l	93
25) 2-Butanone	6.79	43	3499	4.232	ug/l	100
43) Isopropyl Acetate	8.13	43	32212	9.971	ug/l #	90
51) 4-Methyl-2-Pentanone	9.95	43	3511	1.839	ug/l	96

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

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