

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092019\
 Data File : VN058261.D
 Acq On : 20 Sep 2019 19:17
 Operator : JC/SP
 Sample : K4975-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TT101D-20190917

Quant Time: Sep 23 07:04:51 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091819W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 19 09:27:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	475204	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	771348	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	661130	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	263335	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	351621	53.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.26%	
35) Dibromofluoromethane	7.58	113	244697	52.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.36%	
50) Toluene-d8	10.08	98	997255	54.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.16%	
62) 4-Bromofluorobenzene	12.41	95	310139	45.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.40%	
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.84	85	7280	1.865	ug/l	96
9) 1,1,2-Trichlorotrifluoroet	3.74	101	61652	16.048	ug/l	99
12) 1,1-Dichloroethene	3.72	96	11795	3.716	ug/l #	73
16) Acetone	3.80	43	12360	5.502	ug/l	89
27) cis-1,2-Dichloroethene	6.82	96	18311	3.800	ug/l	86
30) Chloroform	7.36	83	4674	0.509	ug/l	93
44) Trichloroethene	8.82	130	424207	102.030	ug/l	100

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

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