

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091119\
 Data File : VN057907.D
 Acq On : 11 Sep 2019 8:27
 Operator : JC/SP
 Sample : K4662-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CL-01-090619

Quant Time: Sep 13 05:44:17 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	305219	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	521138	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	480942	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	249177	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	335594	54.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.58%	
35) Dibromofluoromethane	7.58	113	227102	47.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.44%	
50) Toluene-d8	10.09	98	941226	50.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.80%	
62) 4-Bromofluorobenzene	12.40	95	306418	43.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.12%	
Target Compounds						
16) Acetone	3.80	43	11427	5.639	ug/l	96
18) Methyl Acetate	4.32	43	7240	1.340	ug/l	92
20) Methylene Chloride	4.53	84	48290	11.571	ug/l	95
43) Isopropyl Acetate	8.15	43	81969	7.806	ug/l #	90

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

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