

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN091319\
 Data File : VN057960.D
 Acq On : 12 Sep 2019 11:14
 Operator : JC/SP
 Sample : K4835-03RE
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :

Quant Time: Sep 13 05:09:54 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	247093	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	403142	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	364677	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	193602	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	242738	48.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.00%	
35) Dibromofluoromethane	7.58	113	166610	44.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.56%	
50) Toluene-d8	10.09	98	355533	24.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	49.22%	
62) 4-Bromofluorobenzene	12.41	95	201462	36.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	73.18%	
Target Compounds						
3) Chloromethane	2.03	50	21589	6.199	ug/l	94
11) Tert butyl alcohol	4.77	59	517913	855.029	ug/l	99
16) Acetone	3.80	43	132765	80.923	ug/l	97
18) Methyl Acetate	4.31	43	14144	3.233	ug/l #	83
19) Methyl tert-butyl Ether	5.02	73	1455333	132.911	ug/l	99
25) 2-Butanone	6.83	43	13662	5.753	ug/l	90
30) Chloroform	7.36	83	2594	0.376	ug/l	92

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

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