

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN110918\
 Data File : VN052238.D
 Acq On : 9 Nov 2018 20:15
 Operator : MD\SY
 Sample : J5897-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 CARBON-DNM

Quant Time: Nov 12 09:22:51 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110918W.M
 Quant Title : SW846 8260
 QLast Update : Mon Nov 12 00:34:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	310587	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	512729	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	429815	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	176658	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	194447	45.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.56%	
35) Dibromofluoromethane	7.59	113	154259	46.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.90%	
50) Toluene-d8	10.09	98	630796	48.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.06%	
62) 4-Bromofluorobenzene	12.40	95	191381	45.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.06%	
Target Compounds						
3) Chloromethane	2.06	50	67139	14.242	ug/l	95
18) Methyl Acetate	4.32	43	234783	66.150	ug/l	100
29) Tetrahydrofuran	7.21	42	61374	52.045	ug/l	98

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

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