

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN062218\
 Data File : VN049495.D
 Acq On : 22 Jun 2018 16:49
 Operator : MD\SY
 Sample : J3649-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VOC-DRUM

Quant Time: Jun 22 23:05:22 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061618W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 16 01:02:01 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1561666	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	2437354	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	2036961	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	544574	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	1078179	52.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.10%	
35) Dibromofluoromethane	7.59	113	956746	48.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.04%	
50) Toluene-d8	10.09	98	3405945	45.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.72%	
62) 4-Bromofluorobenzene	12.40	95	834530	33.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	67.22%	
Target Compounds						
3) Chloromethane	2.06	50	18421	0.79	ug/l	95
16) Acetone	3.83	43	2747	0.55	ug/l #	83
18) Methyl Acetate	4.34	43	14548	0.95	ug/l	96
29) Tetrahydrofuran	7.22	42	39381	9.06	ug/l	96

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

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