

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101818\
 Data File : VX005435.D
 Acq On : 18 Oct 2018 17:30
 Operator : JC/MD
 Sample : J5549-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CB-1018-S-TCLP-GRID-50-2

Quant Time: Oct 19 05:00:31 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Oct 12 11:28:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	315930	50.00	ug/l	-0.01
34) 1,4-Difluorobenzene	6.86	114	440670	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	390607	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	190015	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	178160	50.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.62%	
35) Dibromofluoromethane	5.51	113	152734	52.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.66%	
50) Toluene-d8	8.71	98	518273	51.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.44%	
62) 4-Bromofluorobenzene	11.14	95	172420	45.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.62%	
Target Compounds						
16) Acetone	2.45	43	18445	9.098	ug/l	96
18) Methyl Acetate	2.78	43	10060	1.744	ug/l	97
25) 2-Butanone	4.71	43	7277	2.453	ug/l	98
43) Isopropyl Acetate	6.47	43	168306	23.020	ug/l	96

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

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