

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102218\
 Data File : VX005483.D
 Acq On : 22 Oct 2018 21:00
 Operator : JC/MD
 Sample : J5538-24 5X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CNS01

Quant Time: Oct 23 03:14:06 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	5.67	168	247276	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	349518	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	294964	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	144740	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.07	65	155224	56.56	ug/l	0.00
Spiked Amount			Recovery	=	113.12%	
35) Dibromofluoromethane	5.50	113	125051	54.02	ug/l	0.00
Spiked Amount			Recovery	=	108.04%	
50) Toluene-d8	8.71	98	404927	50.95	ug/l	0.00
Spiked Amount			Recovery	=	101.90%	
62) 4-Bromofluorobenzene	11.14	95	132762	44.47	ug/l	0.00
Spiked Amount			Recovery	=	88.94%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.44	43	9749	6.144	ug/l	89
18) Methyl Acetate	2.79	43	6868	1.521	ug/l	98
20) Methylene Chloride	2.86	84	18548	6.894	ug/l	85
43) Isopropyl Acetate	6.47	43	30499	5.259	ug/l	96
51) 4-Methyl-2-Pentanone	8.65	43	79299	19.749	ug/l	94

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

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