

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062518\
 Data File : VX002876.D
 Acq On : 25 Jun 2018 23:01
 Operator : JC/MD
 Sample : J3693-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 W-41

Quant Time: Jun 27 06:56:54 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 25 14:38:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	237977	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	337947	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	312917	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	177652	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	163083	49.58	ug/l	0.00
Spiked Amount			Recovery	=	99.16%	
35) Dibromofluoromethane	5.51	113	131044	48.97	ug/l	0.00
Spiked Amount			Recovery	=	97.94%	
50) Toluene-d8	8.72	98	472088	48.31	ug/l	0.00
Spiked Amount			Recovery	=	96.62%	
62) 4-Bromofluorobenzene	11.14	95	168043	44.59	ug/l	0.00
Spiked Amount			Recovery	=	89.18%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	1.32	50	3522	0.41	ug/l #	60
10) Methyl Iodide	2.53	142	1714	6.79	ug/l #	89
11) Tert butyl alcohol	3.07	59	2476	3.78	ug/l #	92
16) Acetone	2.45	43	3184	2.34	ug/l	95
19) Methyl tert-butyl Ether	3.21	73	9290	1.10	ug/l #	90
65) Chlorobenzene	10.14	112	3031	0.38	ug/l	99

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

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