

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062419\
 Data File : VX010420.D
 Acq On : 24 Jun 2019 18:34
 Operator : JC/SP
 Sample : K3155-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BAT-B10

Quant Time: Jun 25 04:33:20 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 20 03:31:56 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	277405	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	427841	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	388187	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	180237	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	130779	50.28	ug/l	0.00
Spiked Amount			Recovery	=	100.56%	
35) Dibromofluoromethane	5.50	113	129844	49.60	ug/l	0.00
Spiked Amount			Recovery	=	99.20%	
50) Toluene-d8	8.71	98	501709	50.10	ug/l	0.00
Spiked Amount			Recovery	=	100.20%	
62) 4-Bromofluorobenzene	11.13	95	169412	46.86	ug/l	0.00
Spiked Amount			Recovery	=	93.72%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	1.33	50	11584	5.158	ug/l	98
16) Acetone	2.44	43	5429	4.165	ug/l	99
17) Carbon Disulfide	2.57	76	2179	0.338	ug/l #	87
30) Chloroform	5.22	83	2579	0.581	ug/l	98
44) Trichloroethene	7.21	130	4125	1.301	ug/l	98
64) Tetrachloroethene	9.34	164	22126	7.047	ug/l	94

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

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