

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061419\
 Data File : VX010309.D
 Acq On : 14 Jun 2019 18:36
 Operator : JC/SP
 Sample : K3387-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VOC-DRUM

Quant Time: Jun 15 03:02:09 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060719W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 08 02:12:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	305400	50.00	ug/l	-0.02
34) 1,4-Difluorobenzene	6.85	114	454161	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	416144	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	206420	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	132402	44.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.82%	
35) Dibromofluoromethane	5.51	113	120039	42.85	ug/l	0.01
Spiked Amount 50.000			Recovery	=	85.70%	
50) Toluene-d8	8.71	98	533887	50.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.70%	
62) 4-Bromofluorobenzene	11.14	95	189150	48.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.48%	
Target Compounds						
3) Chloromethane	1.33	50	22343	8.006	ug/l	99
17) Carbon Disulfide	2.47	76	21046	2.655	ug/l	97
18) Methyl Acetate	2.82	43	1788929	580.157	ug/l #	84
29) Tetrahydrofuran	5.21	42	110400	84.138	ug/l #	80

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

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