

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX053119\
 Data File : VX009984.D
 Acq On : 31 May 2019 14:50
 Operator : JC/SP
 Sample : K2964-01DL 20X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 COMP-1DL

Quant Time: May 31 15:29:39 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X052819W.M
 Quant Title : SW846 8260
 QLast Update : Tue May 28 13:02:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	167694	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	270237	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	250767	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	118504	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	114343	48.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.54%	
35) Dibromofluoromethane	5.50	113	82947	48.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.64%	
50) Toluene-d8	8.71	98	351559	50.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.92%	
62) 4-Bromofluorobenzene	11.13	95	124644	50.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.60%	
Target Compounds						
16) Acetone	2.45	43	483872	353.539	ug/l	96
18) Methyl Acetate	2.78	43	23105	7.255	ug/l	100
25) 2-Butanone	4.69	43	6808	3.193	ug/l	95
37) Ethyl Acetate	4.84	43	34361	8.994	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX053119\
Data File : VX009984.D
Acq On : 31 May 2019 14:50
Operator : JC/SP
Sample : K2964-01DL 20X
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
COMP-1DL

Quant Time: May 31 15:29:39 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X052819W.M
Quant Title : SW846 8260
QLast Update : Tue May 28 13:02:54 2019
Response via : Initial Calibration

