

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102020\
 Data File : VX018995.D
 Acq On : 20 Oct 2020 23:47
 Operator : JC/SP
 Sample : L4457-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SUTTON-PR-2152-SED

Quant Time: Oct 21 04:19:02 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102020W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 20 19:16:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	162004	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	301370	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	296187	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	145453	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	111953	52.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.14%	
35) Dibromofluoromethane	5.46	113	91401	47.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.98%	
50) Toluene-d8	8.70	98	391480	51.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.16%	
62) 4-Bromofluorobenzene	11.12	95	145182	48.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.94%	
Target Compounds						
16) Acetone	2.42	43	16456	22.193	ug/l	99
20) Methylene Chloride	2.83	84	4427	2.310	ug/l	97
25) 2-Butanone	4.64	43	8196	6.842	ug/l	97
43) Isopropyl Acetate	6.41	43	32119	7.829	ug/l	99
95) Naphthalene	13.82	128	22828	2.118	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102020\
Data File : VX018995.D
Acq On : 20 Oct 2020 23:47
Operator : JC/SP
Sample : L4457-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SUTTON-PR-2152-SED

Quant Time: Oct 21 04:19:02 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102020W.M
Quant Title : SW846 8260
QLast Update : Tue Oct 20 19:16:32 2020
Response via : Initial Calibration

