

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU121318\
 Data File : VU028630.D
 Acq On : 12 Dec 2018 00:29
 Operator : MD/SY
 Sample : J6298-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ENGINE-WASH

Quant Time: Dec 12 05:50:59 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 07 00:39:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	84525	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	153723	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	134004	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	54132	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	69768	54.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.18%	
35) Dibromofluoromethane	4.88	113	51249	52.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.96%	
50) Toluene-d8	7.56	98	197297	51.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.12%	
62) 4-Bromofluorobenzene	10.31	95	65018	45.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.92%	
Target Compounds						
16) Acetone	2.32	43	49725	59.075	ug/l	99
18) Methyl Acetate	2.62	43	5886	2.698	ug/l	94
20) Methylene Chloride	2.70	84	32403	26.114	ug/l	98
25) 2-Butanone	4.27	43	20963	18.224	ug/l	96
43) Isopropyl Acetate	5.55	43	10713	3.115	ug/l #	90
51) 4-Methyl-2-Pentanone	7.46	43	57526	26.518	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU121318\
Data File : VU028630.D
Acq On : 12 Dec 2018 00:29
Operator : MD/SY
Sample : J6298-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ENGINE-WASH

Quant Time: Dec 12 05:50:59 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 07 00:39:42 2018
Response via : Initial Calibration

