

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102621\
 Data File : VU045402.D
 Acq On : 26 Oct 2021 15:05
 Operator : SY/MD
 Sample : M4284-08DL 10X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 20877DL

Quant Time: Oct 26 16:05:26 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 28 12:32:13 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.375	168	216107	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	407390	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	381634	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	175650	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	174110	48.399	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.800%
35) Dibromofluoromethane	5.292	113	126716	49.307	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.620%
50) Toluene-d8	7.899	98	511094	50.333	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.660%
62) 4-Bromofluorobenzene	10.639	95	192714	45.969	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	91.940%
Target Compounds						
					Qvalue	
16) Acetone	2.645	43	790774	370.329	ug/l	98
18) Methyl Acetate	2.954	43	50557	8.101	ug/l #	75
20) Methylene Chloride	3.044	84	1911	0.594	ug/l	87
25) 2-Butanone	4.719	43	83106	27.393	ug/l	99
86) p-Isopropyltoluene	11.796	119	21455	1.858	ug/l	98

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

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