

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040621\
 Data File : VU042949.D
 Acq On : 06 Apr 2021 13:08
 Operator : SY/MD
 Sample : M1903-09DL 2X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 31031DL

Quant Time: Apr 07 01:41:37 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U032421W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 25 05:42:00 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	141530	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	226955	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	210080	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	107223	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	113376	50.25	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.50%	
35) Dibromofluoromethane	5.298	113	79376	46.86	ug/l	0.00
Spiked Amount	50.000		Recovery	=	93.72%	
50) Toluene-d8	7.906	98	283146	47.52	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.04%	
62) 4-Bromofluorobenzene	10.639	95	104060	44.94	ug/l	0.00
Spiked Amount	50.000		Recovery	=	89.88%	
Target Compounds						
						Qvalue
11) Tert butyl alcohol	3.202	59	5985	12.41	ug/l #	88
16) Acetone	2.636	43	504171	447.38	ug/l #	49
18) Methyl Acetate	2.954	43	61225	23.69	ug/l	98
25) 2-Butanone	4.713	43	20635	11.24	ug/l	99

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

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