

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092721\
 Data File : VU044892.D
 Acq On : 27 Sep 2021 16:32
 Operator : SY/MD
 Sample : M3875-15DL 5X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 20882DL

Manual Integrations
 APPROVED

MMDadoda
 9/28/2021 7:04:52 PM

Quant Time: Sep 27 17:25:25 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092421W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 25 09:30:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	200282	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	353604	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	338475	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	167345	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	159826	45.724	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.440%	
35) Dibromofluoromethane	5.292	113	109064	44.960	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.920%	
50) Toluene-d8	7.899	98	449755	49.091	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.180%	
62) 4-Bromofluorobenzene	10.639	95	178641	52.498	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.000%	
Target Compounds					Qvalue	
16) Acetone	2.636	43	117472	59.580	ug/l	99
17) Carbon Disulfide	2.784	76	23977m	3.534	ug/l	
20) Methylene Chloride	3.041	84	164517	51.545	ug/l	97
25) 2-Butanone	4.713	43	24607	8.225	ug/l	99
64) Tetrachloroethene	8.555	164	14372	6.264	ug/l	97

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

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