

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
 Data File : VU045032.D
 Acq On : 30 Sep 2021 07:23
 Operator : SY/MD
 Sample : M3971-11
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 KEARNY-COMP

Quant Time: Sep 30 08:29:56 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	196326	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	355722	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	350264	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	160120	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	159101	48.68	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.36%
35) Dibromofluoromethane	5.292	113	110619	49.30	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.60%
50) Toluene-d8	7.902	98	459123	51.78	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	103.56%
62) 4-Bromofluorobenzene	10.635	95	175753	48.01	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.02%
Target Compounds						
					Qvalue	
16) Acetone	2.629	43	57214	29.49	ug/l	97
18) Methyl Acetate	2.957	43	6397	1.13	ug/l	95
20) Methylene Chloride	3.047	84	9637	3.30	ug/l	95
43) Isopropyl Acetate	5.919	43	14922	2.02	ug/l	98

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

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