

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102221\
 Data File : VU045368.D
 Acq On : 22 Oct 2021 16:35
 Operator : SY/MD
 Sample : M4316-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WC-2-IS

Quant Time: Oct 23 01:13:11 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	212760	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	402470	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	371204	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	169487	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	170410	48.116	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.240%
35) Dibromofluoromethane	5.292	113	122584	48.282	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.560%
50) Toluene-d8	7.899	98	499893	49.831	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.660%
62) 4-Bromofluorobenzene	10.635	95	186401	45.007	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.020%
Target Compounds						
					Qvalue	
16) Acetone	2.636	43	31089	14.788	ug/l	98
18) Methyl Acetate	2.954	43	4310	0.701	ug/l	94
20) Methylene Chloride	3.044	84	16503	5.212	ug/l	95
25) 2-Butanone	4.713	43	7861	2.632	ug/l	95
40) Benzene	5.777	78	13313	1.062	ug/l	99
43) Isopropyl Acetate	5.915	43	29993	3.580	ug/l	99
68) m/p-Xylenes	9.697	106	2996	0.538	ug/l	97
69) o-Xylene	10.105	106	1817	0.329	ug/l	95

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

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