

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102221\
 Data File : VU045369.D
 Acq On : 22 Oct 2021 16:59
 Operator : SY/MD
 Sample : M4316-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WC-2A-ID

Quant Time: Oct 23 01:13:23 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	220177	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	415001	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.424	117	385398	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	171788	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	176323	48.108	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	96.220%		
35) Dibromofluoromethane	5.292	113	127982	48.886	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	97.780%		
50) Toluene-d8	7.899	98	515142	49.801	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	99.600%		
62) 4-Bromofluorobenzene	10.639	95	191476	44.836	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	89.680%		
Target Compounds						Qvalue
16) Acetone	2.642	43	19708	9.059	ug/l	98
18) Methyl Acetate	2.954	43	4965	0.781	ug/l #	87
20) Methylene Chloride	3.041	84	16399	5.005	ug/l	98
25) 2-Butanone	4.723	43	4044	1.308	ug/l	99
40) Benzene	5.771	78	3442	0.266	ug/l	96
43) Isopropyl Acetate	5.919	43	30888	3.575	ug/l	98

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

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