

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092521\
 Data File : VU044846.D
 Acq On : 25 Sep 2021 18:54
 Operator : SY/MD
 Sample : M3829-13
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RE120D1-20210917

Quant Time: Sep 27 01:34:27 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	5.379	168	169768	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	302688	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	282564	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	129271	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	147256	49.700	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.400%	
35) Dibromofluoromethane	5.295	113	99771	48.048	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.100%	
50) Toluene-d8	7.903	98	388560	49.546	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.100%	
62) 4-Bromofluorobenzene	10.636	95	133005	45.661	ug/l	0.00
Spiked Amount	50.000		Recovery	=	91.320%	
Target Compounds						
						Qvalue
9) 1,1,2-Trichlorotrifluo...	2.581	101	22157	10.728	ug/l	96
12) 1,1-Dichloroethene	2.581	96	11166	5.463	ug/l	98
16) Acetone	2.639	43	2673	1.599	ug/l	91
27) cis-1,2-Dichloroethene	4.671	96	5786	2.174	ug/l	89
30) Chloroform	5.092	83	1488	0.325	ug/l	98
44) Trichloroethene	6.549	130	1434202	609.449	ug/l	98
55) 1,1,2-Trichloroethane	8.407	97	1100	0.439	ug/l	91
64) Tetrachloroethene	8.558	164	6678	3.487	ug/l	96

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

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