

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX033021\
 Data File : VX021663.D
 Acq On : 31 Mar 2021 04:58
 Operator : JC/MD
 Sample : M1842-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE104D2-20210323

Quant Time: Mar 31 06:14:02 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 24 12:25:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.629	168	58042	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	104387	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	109009	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	45593	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.035	65	50133	51.91	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 103.82%	
35) Dibromofluoromethane	5.464	113	38705	52.20	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.40%	
50) ToLuene-d8	8.694	98	134746	51.00	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 102.00%	
62) 4-Bromofluorobenzene	11.118	95	51919	50.54	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 101.08%	
Target Compounds						
						Qvalue
16) Acetone	2.423	43	1541	4.22	ug/l #	77
24) 1,1-Dichloroethane	3.676	63	1719	1.38	ug/l #	87
27) cis-1,2-Dichloroethene	4.564	96	29588	39.66	ug/l	96
30) Chloroform	5.176	83	4364	3.32	ug/l	92
42) 1,2-Dichloroethane	6.158	62	2145	1.70	ug/l	85
44) Trichloroethene	7.188	130	157899	206.01	ug/l	93

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

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