

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031722\
 Data File : VX027495.D
 Acq On : 17 Mar 2022 18:35
 Operator : JC/MD
 Sample : N1997-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-101

Quant Time: Mar 18 02:06:19 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 17 15:02:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	335066	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	545850	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	490983	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	224014	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	204593	49.566	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	99.140%
35) Dibromofluoromethane	5.391	113	183545	52.264	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	104.520%
50) Toluene-d8	8.653	98	638901	48.076	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.160%
62) 4-Bromofluorobenzene	11.085	95	231744	47.444	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	94.880%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.965	59	15198	19.320	ug/l	99
16) Acetone	2.380	43	4277	2.542	ug/l	97
19) Methyl tert-butyl Ether	3.117	73	69492	6.620	ug/l	100
30) Chloroform	5.105	83	7642	1.198	ug/l	97
40) Benzene	6.044	78	33913	2.463	ug/l	96
42) 1,2-Dichloroethane	6.092	62	15122	2.931	ug/l	95
68) m/p-Xylenes	10.305	106	3351	0.492	ug/l	98

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

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