

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031722\
 Data File : VX027506.D
 Acq On : 17 Mar 2022 22:55
 Operator : JC/MD
 Sample : N1997-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-112

Quant Time: Mar 18 02:08:19 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	329015	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	544201	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	489557	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	222589	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	205150	50.615	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 101.240%	
35) Dibromofluoromethane	5.391	113	179275	51.203	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 102.400%	
50) Toluene-d8	8.653	98	638111	48.162	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 96.320%	
62) 4-Bromofluorobenzene	11.085	95	227109	46.636	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 93.280%	
Target Compounds						Qvalue
11) Tert butyl alcohol	2.965	59	220139	284.988	ug/l #	72
16) Acetone	2.380	43	6358	3.848	ug/l #	84
19) Methyl tert-butyl Ether	3.117	73	13154794	1276.195	ug/l	92
22) Diisopropyl ether	3.763	45	160855	17.229	ug/l	91
30) Chloroform	5.111	83	2526	0.403	ug/l	95
31) Cyclohexane	5.483	56	9357	1.859	ug/l #	81
39) Methylcyclohexane	7.379	83	2495	0.415	ug/l #	81
40) Benzene	6.043	78	343476	25.022	ug/l	99

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

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