

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031722\
 Data File : VX027503.D
 Acq On : 17 Mar 2022 21:44
 Operator : JC/MD
 Sample : N1997-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-107

Quant Time: Mar 18 02:07:46 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	332545	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	549330	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	489591	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	214143	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	207503	50.652	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	101.300%
35) Dibromofluoromethane	5.391	113	181660	51.399	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	102.800%
50) Toluene-d8	8.653	98	642125	48.012	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.020%
62) 4-Bromofluorobenzene	11.085	95	224514	45.672	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	91.340%
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	5444	3.260	ug/l	92
19) Methyl tert-butyl Ether	3.123	73	12300	1.181	ug/l	91
30) Chloroform	5.098	83	4948	0.782	ug/l	94
40) Benzene	6.043	78	5112	0.369	ug/l	99
42) 1,2-Dichloroethane	6.098	62	8337	1.605	ug/l	98

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

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