

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
 Data File : VU049539.D
 Acq On : 30 Jun 2022 02:45
 Operator : SY/MD
 Sample : N3522-02
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-TT-MW149I1-330-332

Quant Time: Jun 30 04:14:06 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 28 10:35:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.372	168	243962	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.250	114	428354	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	390071	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	174524	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	172397	45.885	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	91.760%
35) Dibromofluoromethane	5.288	113	137308	45.817	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.640%
50) Toluene-d8	7.899	98	516835	48.352	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.700%
62) 4-Bromofluorobenzene	10.632	95	184128	44.968	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	89.940%
Target Compounds						
					Qvalue	
16) Acetone	2.636	43	8640	4.560	ug/l	98
17) Carbon Disulfide	2.793	76	3973	0.481	ug/l #	87
19) Methyl tert-butyl Ether	3.366	73	19259	1.954	ug/l	99
60) Dibromochloromethane	8.809	129	1703	0.508	ug/l	98

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

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