

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062922\
 Data File : VU049520.D
 Acq On : 29 Jun 2022 17:54
 Operator : SY/MD
 Sample : N3522-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-TT-MW14911-350-352

Quant Time: Jun 30 00:49:30 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	262637	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	461968	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	422529	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	194717	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	182554	45.133	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.260%
35) Dibromofluoromethane	5.288	113	144342	44.660	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.320%
50) Toluene-d8	7.896	98	563434	48.876	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.760%
62) 4-Bromofluorobenzene	10.632	95	206775	46.825	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.640%
Target Compounds						
					Qvalue	
16) Acetone	2.626	43	26197	12.843	ug/l	98
17) Carbon Disulfide	2.790	76	3009	0.338	ug/l #	89
25) 2-Butanone	4.716	43	8417	2.753	ug/l	91
40) Benzene	5.777	78	8358	0.591	ug/l	99
52) Toluene	7.970	92	22836	2.715	ug/l	100
67) Ethyl Benzene	9.571	91	7423	0.468	ug/l	97
68) m/p-Xylenes	9.693	106	9664	1.647	ug/l	93
69) o-Xylene	10.102	106	3647	0.626	ug/l	89
84) 1,2,4-Trimethylbenzene	11.468	105	16494	1.386	ug/l	72
95) Naphthalene	14.089	128	39813	4.257	ug/l	99

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

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