

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062922\
 Data File : VX029868.D
 Acq On : 29 Jun 2022 17:40
 Operator : JC/MD
 Sample : N3485-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT205S1-20220623

Quant Time: Jun 29 23:00:51 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 05:54:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	217122	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	374275	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	313300	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	131652	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	119152	43.353	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.700%
35) Dibromofluoromethane	5.391	113	105327	45.030	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.060%
50) Toluene-d8	8.653	98	422847	48.189	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.380%
62) 4-Bromofluorobenzene	11.085	95	138461	44.640	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	89.280%
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.325	96	776	0.347	ug/l #	63
16) Acetone	2.386	43	2453	1.824	ug/l	91
19) Methyl tert-butyl Ether	3.117	73	1929	0.256	ug/l	99
24) 1,1-Dichloroethane	3.617	63	1743	0.396	ug/l #	91
30) Chloroform	5.086	83	1902	0.441	ug/l	99
44) Trichloroethene	7.129	130	16249	5.998	ug/l	99
64) Tetrachloroethene	9.281	164	988	0.523	ug/l	89

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

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