

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062922\
 Data File : VX029867.D
 Acq On : 29 Jun 2022 17:17
 Operator : JC/MD
 Sample : N3485-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT161S1-20220623

Quant Time: Jun 29 23:00:40 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	245045	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	422471	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	353620	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	151124	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	132682	42.775	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	85.560%
35) Dibromofluoromethane	5.391	113	118126	44.741	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.480%
50) Toluene-d8	8.653	98	476802	48.139	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.280%
62) 4-Bromofluorobenzene	11.085	95	157701	45.042	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.080%
Target Compounds						
					Qvalue	
3) Chloromethane	1.294	50	615	0.223	ug/l #	79
12) 1,1-Dichloroethene	2.319	96	1484	0.588	ug/l #	75
19) Methyl tert-butyl Ether	3.123	73	4597	0.540	ug/l #	87
24) 1,1-Dichloroethane	3.617	63	12932	2.605	ug/l	97
30) Chloroform	5.092	83	1588	0.326	ug/l	86
44) Trichloroethene	7.129	130	9891	3.235	ug/l	94
64) Tetrachloroethene	9.275	164	902	0.423	ug/l #	70

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

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