

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062222\
 Data File : VX029704.D
 Acq On : 22 Jun 2022 17:10
 Operator : JC/MD
 Sample : N3293-19
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 30986

Quant Time: Jun 22 17:34:53 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 20 01:31:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.555	168	267190	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	450706	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	388662	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	176789	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	149733	42.335	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	84.660%
35) Dibromofluoromethane	5.391	113	133138	45.328	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.660%
50) Toluene-d8	8.652	98	507915	47.335	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.680%
62) 4-Bromofluorobenzene	11.079	95	183906	45.449	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.900%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.964	59	40369	44.780	ug/l	97
16) Acetone	2.379	43	206629	146.570	ug/l	100
17) Carbon Disulfide	2.526	76	39880	5.683	ug/l	97
18) Methyl Acetate	2.702	43	334091	89.633	ug/l	97
20) Methylene Chloride	2.788	84	3564	1.135	ug/l	97
25) 2-Butanone	4.556	43	73186	32.843	ug/l	96
29) Tetrahydrofuran	5.019	42	5210	3.508	ug/l	87
40) Benzene	6.037	78	29212	2.479	ug/l	95
52) Toluene	8.720	92	12928	1.713	ug/l	100
59) 2-Hexanone	9.439	43	4184	1.353	ug/l	93
68) m/p-Xylenes	10.305	106	5825	1.089	ug/l	98
69) o-Xylene	10.646	106	5421	1.026	ug/l	92

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

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