

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091422\
 Data File : VX031342.D
 Acq On : 15 Sep 2022 06:46
 Operator : JC/MD
 Sample : N4579-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CHERRY-HILL-COMP-11

Quant Time: Sep 15 11:12:24 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091422W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 15 01:50:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	55533	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	103352	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	98099	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	41780	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	56709	41.604	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	83.200%
35) Dibromofluoromethane	5.385	113	41295	37.171	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	74.340%#
50) Toluene-d8	8.653	98	162954	39.102	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	78.200%#
62) 4-Bromofluorobenzene	11.079	95	55509	36.374	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	72.740%#
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	17081	29.354	ug/l	97
20) Methylene Chloride	2.794	84	3369	1.369	ug/l	92
37) Ethyl Acetate	4.715	43	10928	5.469	ug/l	98
43) Isopropyl Acetate	6.342	43	52266	15.323	ug/l	97

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

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