

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091422\
 Data File : VX031344.D
 Acq On : 15 Sep 2022 07:32
 Operator : JC/MD
 Sample : N4579-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CHERRY-HILL-COMP-13

Quant Time: Sep 15 11:12:58 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	56682	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	104123	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	100339	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	42294	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	59274	42.604	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	85.200%		
35) Dibromofluoromethane	5.391	113	43834	39.164	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	78.320%		
50) Toluene-d8	8.653	98	173887	41.417	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	82.840%#		
62) 4-Bromofluorobenzene	11.079	95	59632	38.787	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	77.580%#		
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	23367	39.342	ug/l	98
20) Methylene Chloride	2.788	84	5065	3.203	ug/l	98
25) 2-Butanone	4.574	43	3158	3.223	ug/l	97
37) Ethyl Acetate	4.715	43	10926	5.427	ug/l	97
43) Isopropyl Acetate	6.342	43	47942	13.951	ug/l	96

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

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