

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091422\
 Data File : VX031343.D
 Acq On : 15 Sep 2022 07:09
 Operator : JC/MD
 Sample : N4579-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CHERRY-HILL-COMP-12

Quant Time: Sep 15 11:12:45 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	54632	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	100786	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	95706	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	39473	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	56441	42.090	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	84.180%		
35) Dibromofluoromethane	5.391	113	41667	38.460	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	76.920%		
50) Toluene-d8	8.653	98	163270	40.176	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	80.360%#		
62) 4-Bromofluorobenzene	11.079	95	55321	37.174	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	74.340%#		
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	23790	41.557	ug/l	93
20) Methylene Chloride	2.794	84	4827	3.139	ug/l	93
25) 2-Butanone	4.581	43	3248	3.440	ug/l	93
37) Ethyl Acetate	4.721	43	11476	5.889	ug/l #	90
43) Isopropyl Acetate	6.342	43	49571	14.903	ug/l	96

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

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