

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091422\
 Data File : VX031346.D
 Acq On : 15 Sep 2022 08:19
 Operator : JC/MD
 Sample : N4579-20
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CHERRY-HILL-COMP-15

Quant Time: Sep 15 11:13:20 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	54191	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	99644	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	95983	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	39949	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	57727	43.399	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.800%
35) Dibromofluoromethane	5.391	113	41590	38.829	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	77.660%
50) Toluene-d8	8.653	98	164508	40.944	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	81.880%#
62) 4-Bromofluorobenzene	11.079	95	55024	37.398	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	74.800%#
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	7319	12.889	ug/l	98
20) Methylene Chloride	2.794	84	5121	3.531	ug/l	94
37) Ethyl Acetate	4.715	43	11459	5.948	ug/l	94
43) Isopropyl Acetate	6.342	43	49090	14.927	ug/l	96

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

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