

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091422\
 Data File : VX031333.D
 Acq On : 15 Sep 2022 02:06
 Operator : JC/MD
 Sample : N4578-37
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CHERRY-HILL-COMP-9

Quant Time: Sep 15 06:10:13 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	57973	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	106116	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	103110	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	42331	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	60694	42.653	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	85.300%
35) Dibromofluoromethane	5.391	113	44118	38.677	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	77.360%
50) Toluene-d8	8.653	98	173595	40.571	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	81.140%#
62) 4-Bromofluorobenzene	11.079	95	59192	37.777	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	75.560%#
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	14451	23.789	ug/l	98
20) Methylene Chloride	2.788	84	3518	1.370	ug/l	97
37) Ethyl Acetate	4.720	43	12629	6.155	ug/l	97
43) Isopropyl Acetate	6.342	43	56513	16.137	ug/l	97

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

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