

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091522\
 Data File : VN074470.D
 Acq On : 15 Sep 2022 18:47
 Operator : JC\MD
 Sample : N4578-12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CHERRY-HILL-COMP-3

Quant Time: Sep 15 19:26:05 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091522W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 15 19:20:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	440344	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	854900	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	867178	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	379673	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	321311	43.005	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.000%
35) Dibromofluoromethane	7.951	113	264219	46.526	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.060%
50) Toluene-d8	10.380	98	992662	45.198	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	90.400%
62) 4-Bromofluorobenzene	12.674	95	404737	53.806	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	107.620%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	115575	24.174	ug/l	96
20) Methylene Chloride	5.010	84	24643	3.405	ug/l #	86
37) Ethyl Acetate	7.333	43	119617	7.706	ug/l	97
43) Isopropyl Acetate	8.492	43	591620	24.330	ug/l #	80
68) m/p-Xylenes	11.904	106	4960	0.336	ug/l	100

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

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