

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091522\
 Data File : VN074468.D
 Acq On : 15 Sep 2022 17:58
 Operator : JC\MD
 Sample : N4578-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CHERRY-HILL-COMP-1

Quant Time: Sep 15 19:30:28 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	452133	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	877137	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	896481	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	389611	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	329217	42.914	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	85.820%		
35) Dibromofluoromethane	7.951	113	277337	47.598	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.200%		
50) Toluene-d8	10.381	98	1029084	45.668	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	91.340%		
62) 4-Bromofluorobenzene	12.669	95	411390	53.304	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	106.600%		
Target Compounds						
16) Acetone	4.228	43	125140	25.492	ug/l	90
20) Methylene Chloride	5.010	84	27002	3.631	ug/l	90
37) Ethyl Acetate	7.340	43	119984	7.533	ug/l	97
43) Isopropyl Acetate	8.498	43	547205	21.933	ug/l #	84
68) m/p-Xylenes	11.904	106	4041	0.265	ug/l	84

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

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