

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092521\
 Data File : VN068674.D
 Acq On : 25 Sep 2021 18:00
 Operator : JC/MD
 Sample : PB139363ZHE#02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB139363ZHE#02

Quant Time: Sep 27 05:58:20 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N092321W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 23 13:52:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.091	168	11013	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.971	114	19264	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	16502	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.680	152	7120	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	8447	49.132	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	98.260%		
35) Dibromofluoromethane	8.024	113	6061	46.946	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	93.900%		
50) Toluene-d8	10.443	98	23925	43.791	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	87.580%		
62) 4-Bromofluorobenzene	12.736	95	6562	35.119	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	70.240%		
Target Compounds						Qvalue
16) Acetone	4.293	43	105107	1754.644	ug/l	93
18) Methyl Acetate	4.875	43	10508	41.553	ug/l #	71
20) Methylene Chloride	5.103	84	30175	239.144	ug/l	96
25) 2-Butanone	7.343	43	25078	239.266	ug/l	97
43) Isopropyl Acetate	8.568	43	66300	175.462	ug/l #	83
51) 4-Methyl-2-Pentanone	10.333	43	13685	59.887	ug/l	90
80) 1,3,5-Trimethylbenzene	13.061	105	1797	3.584	ug/l	92

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

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