

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122923\
 Data File : VN080614.D
 Acq On : 29 Dec 2023 14:47
 Operator : JC\MD
 Sample : PB158132ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB158132ZHE#06

Quant Time: Dec 29 23:43:20 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122723W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 28 03:07:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	277633	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	579560	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	695008	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	279354	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	240080	67.414	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 134.820%#		
35) Dibromofluoromethane	8.165	113	220857	58.921	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 117.840%		
50) Toluene-d8	10.571	98	856576	65.450	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 130.900%#		
62) 4-Bromofluorobenzene	12.853	95	310414	70.476	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 140.960%#		
Target Compounds						
						Qvalue
3) Chloromethane	2.365	50	9436	1.802	ug/l	97
16) Acetone	4.424	43	56450	58.312	ug/l	95
20) Methylene Chloride	5.283	84	27907	9.035	ug/l #	85
25) 2-Butanone	7.483	43	14744	8.149	ug/l #	76
29) Tetrahydrofuran	7.847	42	7700	6.525	ug/l #	84
37) Ethyl Acetate	7.565	43	6573	1.510	ug/l #	91
43) Isopropyl Acetate	8.688	43	221838	31.442	ug/l #	85
51) 4-Methyl-2-Pentanone	10.447	43	25462	6.345	ug/l #	80

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

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