

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060925\
 Data File : VN086909.D
 Acq On : 09 Jun 2025 16:34
 Operator : JC\MD
 Sample : Q2228-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP08-MHI-WC

Quant Time: Jun 10 03:35:57 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 07 02:12:50 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	329169	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	589208	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	503460	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	243469	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	192160	43.601	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	87.200%	
35) Dibromofluoromethane	8.177	113	166924	47.805	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	95.600%	
50) Toluene-d8	10.570	98	704465	50.963	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	101.920%	
62) 4-Bromofluorobenzene	12.847	95	249195	48.522	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	97.040%	
Target Compounds						
					Qvalue	
8) Diethyl Ether	3.989	74	81216	32.634	ug/l	91
16) Acetone	4.447	43	28790	12.318	ug/l	96
18) Methyl Acetate	5.047	43	29734	4.366	ug/l	97
20) Methylene Chloride	5.300	84	7293	1.667	ug/l	93
37) Ethyl Acetate	7.571	43	20490	3.094	ug/l	98
43) Isopropyl Acetate	8.700	43	12867	1.210	ug/l #	95

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

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