

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081325\
 Data File : VN087536.D
 Acq On : 13 Aug 2025 15:36
 Operator : JC\MD
 Sample : Q2808-04RE
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-7RE

Quant Time: Aug 14 04:02:02 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071625W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 17 02:56:13 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.212	168	221491	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.088	114	499677	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.847	117	478898	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	229716	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.565	65	243377	64.758	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 129.520%#	
35) Dibromofluoromethane	8.153	113	174728	50.693	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.380%	
50) Toluene-d8	10.547	98	655545	53.318	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 106.640%	
62) 4-Bromofluorobenzene	12.829	95	245693	54.088	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 108.180%	
Target Compounds						
					Qvalue	
16) Acetone	4.424	43	86221	48.930	ug/l	100
18) Methyl Acetate	5.024	43	53471	12.078	ug/l	97
20) Methylene Chloride	5.253	84	21039	6.506	ug/l	88
25) 2-Butanone	7.477	43	23709	8.708	ug/l	92
37) Ethyl Acetate	7.559	43	34715	5.278	ug/l	100
43) Isopropyl Acetate	8.688	43	23495	2.301	ug/l #	88

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

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