



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 2VRE / pdb_00002vre
Title : crystal structure of human peroxisomal delta3,5,delta2,4-dienoyl coa isomerase
Authors : Yue, W.; Guo, K.; von Delft, F.; Pilka, E.; Murray, J.; Roos, A.; Kochan, G.; Bountra, C.; Arrowsmith, C.; Wikstrom, M.; Edwards, A.; Oppermann, U.
Deposited on : 2008-03-31
Resolution : 1.95 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

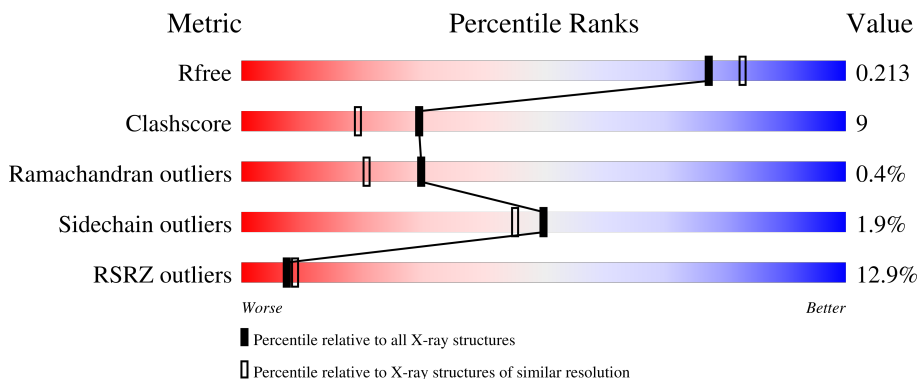
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	296	<div style="display: flex; align-items: center;"> <div style="width: 10%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 81%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 8%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">81% 11% 8%</p>
1	B	296	<div style="display: flex; align-items: center;"> <div style="width: 12%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 73%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 14%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">73% 14% • 12%</p>
1	C	296	<div style="display: flex; align-items: center;"> <div style="width: 11%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 78%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">78% 10% 12%</p>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6452 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called DELTA(3,5)-DELTA(2,4)-DIENOYL-COA ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	272	2102	1321	368	397	16	0	7	0
1	B	260	1968	1236	345	371	16	0	2	0
1	C	261	1983	1241	354	373	15	0	1	0

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cl	0	0
			2	2		
2	B	2	Total	Cl	0	0
			2	2		
2	C	2	Total	Cl	0	0
			2	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	166	Total	O	0	0
			166	166		
3	B	128	Total	O	0	0
			128	128		
3	C	99	Total	O	0	0
			99	99		

THR
THR
GLU
ASN
LYS
GLU
LEU
LYS
THR

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	127.58Å 137.25Å 96.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	93.25 – 1.95 93.25 – 1.95	Depositor EDS
% Data completeness (in resolution range)	100.0 (93.25-1.95) 100.0 (93.25-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 1.88Å)	Xtrriage
Refinement program	REFMAC 5.4.0066	Depositor
R, R_{free}	0.167 , 0.202 0.185 , 0.213	Depositor DCC
R_{free} test set	3442 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	31.6	Xtrriage
Anisotropy	0.289	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6452	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.91	2/2154 (0.1%)	0.94	0/2919
1	B	0.87	2/2003 (0.1%)	0.94	3/2714 (0.1%)
1	C	0.78	0/2018	0.94	4/2735 (0.1%)
All	All	0.86	4/6175 (0.1%)	0.94	7/8368 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	232	LEU	N-CA	7.71	1.55	1.46
1	A	21	GLN	CD-OE1	5.47	1.33	1.23
1	B	50	ASN	CG-OD1	5.26	1.33	1.23
1	B	40	HIS	ND1-CE1	5.21	1.37	1.32

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	40	HIS	CB-CG-CD2	-6.34	122.96	131.20
1	C	47	ASN	N-CA-C	6.15	117.81	110.44
1	B	171	GLU	N-CA-C	5.70	118.23	111.33
1	B	40	HIS	CB-CG-ND1	5.48	130.92	122.70
1	C	208	ALA	N-CA-C	5.38	117.14	111.28
1	C	22	SER	N-CA-CB	-5.13	101.78	110.50
1	C	22	SER	N-CA-C	5.10	125.28	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2102	0	2081	37	0
1	B	1968	0	1940	44	0
1	C	1983	0	1958	31	0
2	A	2	0	0	0	1
2	B	2	0	0	0	1
2	C	2	0	0	0	0
3	A	166	0	0	13	0
3	B	128	0	0	14	0
3	C	99	0	0	8	0
All	All	6452	0	5979	103	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (103) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157[B]:ASP:OD2	1:B:201:PHE:CE2	1.65	1.48
3:A:2097:HOH:O	1:C:280:ASP:HB2	1.22	1.30
1:B:25:PRO:CD	3:B:2002:HOH:O	1.65	1.29
1:A:157[B]:ASP:OD2	1:B:201:PHE:HE2	0.89	1.21
1:A:131[B]:GLU:CD	3:A:2074:HOH:O	1.81	1.18
1:C:67:PHE:CE1	1:C:130:ILE:HG12	1.77	1.18
1:A:33[B]:ARG:NH2	1:A:45:GLN:OE1	1.79	1.14
1:B:101:ILE:HA	3:B:2034:HOH:O	1.53	1.09
1:C:74:ALA:HA	3:C:2016:HOH:O	1.51	1.06
1:A:131[B]:GLU:OE1	3:A:2074:HOH:O	1.71	1.04
1:B:25:PRO:HD2	3:B:2002:HOH:O	1.36	1.03
1:A:157[B]:ASP:OD1	1:B:197:ASN:ND2	1.92	1.02
1:C:63:MET:SD	1:C:67:PHE:CE2	2.51	1.02
1:B:24:ALA:HA	3:B:2002:HOH:O	1.64	0.97
1:B:25:PRO:HD3	3:B:2002:HOH:O	1.35	0.97
1:C:63:MET:O	1:C:67:PHE:HD2	1.47	0.97
1:B:219:ARG:HD2	3:C:2058:HOH:O	1.66	0.94
1:A:157[B]:ASP:OD1	1:B:197:ASN:CB	2.16	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:MET:O	1:C:67:PHE:CD2	2.25	0.89
1:A:157[B]:ASP:OD2	1:B:201:PHE:CZ	2.27	0.87
1:C:198:GLU:OE1	3:C:2059:HOH:O	1.93	0.85
1:A:157[B]:ASP:OD1	1:B:197:ASN:CG	2.21	0.83
1:A:37:ALA:O	1:A:38:GLN:HB2	1.78	0.83
1:C:39:LYS:HE2	1:C:40:HIS:NE2	1.94	0.83
1:C:63:MET:SD	1:C:67:PHE:HE2	2.00	0.83
1:C:67:PHE:CE1	1:C:130:ILE:CG1	2.60	0.82
1:C:224:LYS:HE2	3:C:2020:HOH:O	1.82	0.80
1:B:163:GLN:HB3	3:B:2102:HOH:O	1.80	0.80
1:A:38:GLN:HG2	1:A:39:LYS:H	1.53	0.74
1:A:229:ASP:OD1	3:A:2127:HOH:O	2.05	0.73
1:A:157[B]:ASP:OD1	1:B:197:ASN:HB3	1.89	0.72
1:B:85:GLY:HA3	3:B:2029:HOH:O	1.90	0.71
1:A:157[B]:ASP:CG	1:B:201:PHE:HE2	1.95	0.71
1:A:45:GLN:CD	3:A:2021:HOH:O	2.33	0.70
1:A:188:LYS:CE	3:A:2102:HOH:O	2.41	0.69
1:C:63:MET:SD	1:C:67:PHE:CZ	2.86	0.68
1:C:76:CYS:O	1:C:135:LYS:HE2	1.95	0.66
1:A:188:LYS:HE3	3:A:2102:HOH:O	1.98	0.64
1:A:37:ALA:O	1:A:38:GLN:CB	2.47	0.62
1:A:255:TYR:OH	3:A:2144:HOH:O	2.14	0.62
1:C:172:VAL:HG13	3:C:2061:HOH:O	1.98	0.62
1:A:188:LYS:HE2	3:A:2102:HOH:O	2.00	0.60
1:B:127:PHE:CE2	1:B:180:VAL:HG11	2.36	0.60
1:A:33[B]:ARG:NH1	1:A:35:THR:HG21	2.17	0.60
1:B:127:PHE:HE2	1:B:180:VAL:HG11	1.68	0.59
1:A:131[B]:GLU:CD	3:A:2075:HOH:O	2.47	0.57
1:B:36:SER:HB3	1:B:42:LEU:HD23	1.85	0.57
1:A:45:GLN:HE21	1:A:84:ALA:CA	2.19	0.56
1:B:100:ASP:HA	3:B:2042:HOH:O	2.06	0.55
1:A:157[B]:ASP:CG	1:B:197:ASN:HB3	2.32	0.55
1:A:125:GLU:OE2	1:A:132:ARG:NH1	2.25	0.54
1:C:67:PHE:CZ	1:C:130:ILE:HG12	2.38	0.54
1:C:142:HIS:CE1	1:C:224:LYS:HE3	2.43	0.54
1:C:67:PHE:HE1	1:C:130:ILE:HG12	1.58	0.54
1:B:127:PHE:HD2	1:B:180:VAL:HG12	1.72	0.54
1:C:39:LYS:HE2	1:C:40:HIS:CE1	2.43	0.54
1:A:263:GLU:OE1	3:A:2155:HOH:O	2.18	0.54
1:C:67:PHE:CD2	1:C:129:VAL:HG23	2.43	0.54
1:C:87:MET:HE3	1:C:144:GLY:HA3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:PRO:HB3	3:B:2034:HOH:O	2.08	0.53
1:C:67:PHE:CZ	1:C:130:ILE:CG1	2.91	0.52
1:A:45:GLN:HE21	1:A:84:ALA:HA	1.75	0.52
1:B:58:VAL:HG23	3:B:2015:HOH:O	2.10	0.52
1:A:63:MET:HB2	1:A:126:THR:HG23	1.91	0.52
1:C:115:TYR:CZ	1:C:119:ILE:HD11	2.45	0.52
1:C:125:GLU:OE2	1:C:132:ARG:NH1	2.43	0.51
1:B:280:ASP:OD1	1:B:280:ASP:N	2.44	0.50
1:A:33[B]:ARG:NH1	1:A:35:THR:CG2	2.75	0.49
1:B:108:ASP:HB2	3:B:2039:HOH:O	2.11	0.49
1:A:58:VAL:HG22	1:A:61:ARG:NH2	2.28	0.49
1:A:152:LEU:C	1:A:152:LEU:HD23	2.38	0.48
1:B:127:PHE:CD2	1:B:180:VAL:CG1	2.96	0.48
1:C:67:PHE:CG	1:C:129:VAL:HG23	2.49	0.48
1:B:100:ASP:CA	3:B:2042:HOH:O	2.60	0.47
1:A:115:TYR:CZ	1:A:119:ILE:HD11	2.50	0.46
1:B:218:SER:HB2	3:C:2059:HOH:O	2.16	0.46
1:B:152:LEU:C	1:B:152:LEU:HD23	2.41	0.45
1:B:85:GLY:CA	3:B:2029:HOH:O	2.55	0.45
1:A:87:MET:HE3	1:A:144:GLY:HA3	1.98	0.45
1:B:63:MET:HB2	1:B:126:THR:HG23	1.97	0.45
1:B:127:PHE:CD2	1:B:180:VAL:HG12	2.51	0.45
1:C:63:MET:C	1:C:67:PHE:HD2	2.22	0.45
1:C:67:PHE:CD2	1:C:129:VAL:CG2	3.00	0.44
1:A:172[A]:VAL:HG13	3:A:2092:HOH:O	2.18	0.44
1:B:86:LYS:N	3:B:2029:HOH:O	2.41	0.44
1:A:229:ASP:CG	3:A:2127:HOH:O	2.56	0.44
1:A:45:GLN:NE2	1:A:84:ALA:HB2	2.33	0.44
1:B:70:ILE:O	1:B:135:LYS:HE2	2.18	0.43
1:B:174:VAL:HG12	1:B:174:VAL:O	2.17	0.43
1:B:206:MET:HG3	1:B:210:GLU:HB3	1.98	0.43
1:B:45:GLN:NE2	1:B:84:ALA:HB2	2.34	0.42
1:C:224:LYS:CE	3:C:2020:HOH:O	2.54	0.42
1:B:67:PHE:CZ	1:B:81:ILE:HD11	2.55	0.42
1:B:25:PRO:O	1:B:26:ASP:CB	2.68	0.42
1:C:172:VAL:HG22	3:C:2061:HOH:O	2.20	0.42
1:C:280:ASP:OD1	1:C:280:ASP:N	2.53	0.41
1:B:39:LYS:HG2	1:B:40:HIS:CD2	2.56	0.41
1:C:39:LYS:HG2	1:C:40:HIS:CD2	2.54	0.41
1:B:279:GLN:C	1:B:281:LEU:H	2.29	0.41
1:B:80:VAL:HG21	1:B:232:LEU:HD23	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:VAL:HG21	1:C:232:LEU:HD23	2.02	0.40
1:B:115:TYR:CZ	1:B:119:ILE:HD11	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1291:CL:CL	2:B:1284:CL:CL[4_555]	0.29	1.91

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	277/296 (94%)	272 (98%)	3 (1%)	2 (1%)	18	10
1	B	259/296 (88%)	250 (96%)	9 (4%)	0	100	100
1	C	260/296 (88%)	254 (98%)	5 (2%)	1 (0%)	30	21
All	All	796/888 (90%)	776 (98%)	17 (2%)	3 (0%)	30	21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	C	107	ASP
1	A	107	ASP

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	225/251 (90%)	219 (97%)	6 (3%)	39	32
1	B	208/251 (83%)	206 (99%)	2 (1%)	68	67
1	C	209/251 (83%)	203 (97%)	6 (3%)	37	29
All	All	642/753 (85%)	628 (98%)	14 (2%)	50	40

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	28[A]	SER
1	A	28[B]	SER
1	A	135	LYS
1	A	218	SER
1	A	281	LEU
1	B	228	LEU
1	B	280	ASP
1	C	58	VAL
1	C	61[A]	ARG
1	C	61[B]	ARG
1	C	218	SER
1	C	280	ASP
1	C	282	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	168	GLN
1	A	193	GLN
1	A	277	GLN
1	B	45	GLN
1	B	163	GLN
1	C	142	HIS
1	C	252	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	272/296 (91%)	1.07	31 (11%) 10 11	21, 36, 54, 78	8 (2%)
1	B	260/296 (87%)	1.18	37 (14%) 6 7	21, 36, 54, 64	5 (1%)
1	C	261/296 (88%)	1.11	34 (13%) 7 8	28, 36, 52, 57	1 (0%)
All	All	793/888 (89%)	1.12	102 (12%) 7 9	21, 36, 53, 78	14 (1%)

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	67	PHE	9.5
1	B	101	ILE	8.1
1	B	105	LYS	6.6
1	C	25	PRO	6.1
1	A	287	ALA	5.6
1	A	289	THR	5.3
1	A	38	GLN	5.1
1	B	97	MET	5.0
1	A	283	LYS	4.9
1	C	24	ALA	4.9
1	A	288	THR	4.7
1	B	98	ALA	4.5
1	A	101	ILE	4.4
1	B	107	ASP	4.4
1	C	37	ALA	4.4
1	B	99	SER	4.3
1	B	104	PRO	4.3
1	B	102	LEU	4.1
1	C	107	ASP	4.0
1	B	282	VAL	4.0
1	C	23	MET	4.0
1	B	37	ALA	3.9
1	C	102	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	106	GLY	3.9
1	A	285	VAL	3.9
1	B	280	ASP	3.9
1	A	19	TYR	3.8
1	A	131[A]	GLU	3.8
1	C	280	ASP	3.7
1	C	261	VAL	3.7
1	C	94	LEU	3.6
1	B	61	ARG	3.6
1	C	282	VAL	3.6
1	A	18	LEU	3.5
1	C	97	MET	3.4
1	B	106	GLY	3.3
1	B	180	VAL	3.3
1	C	104	PRO	3.2
1	C	103	GLN	3.2
1	A	37	ALA	3.2
1	B	255	TYR	3.1
1	A	279	GLN	3.1
1	B	100	ASP	3.1
1	C	260	SER	3.1
1	B	23	MET	3.0
1	C	61[A]	ARG	3.0
1	A	78	ALA	3.0
1	B	103	GLN	2.9
1	A	106	GLY	2.9
1	A	282	VAL	2.8
1	B	109	VAL	2.8
1	C	105	LYS	2.8
1	B	78	ALA	2.8
1	A	102	LEU	2.8
1	B	114	TRP	2.8
1	A	41	VAL	2.8
1	C	101	ILE	2.8
1	A	174	VAL	2.8
1	C	26	ASP	2.7
1	A	262	ALA	2.7
1	C	160	TYR	2.7
1	B	28	SER	2.7
1	B	261	VAL	2.7
1	B	123	TYR	2.6
1	A	286	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	263	GLU	2.6
1	A	40	HIS	2.6
1	C	172	VAL	2.6
1	C	88	PHE	2.5
1	C	279	GLN	2.5
1	C	256	SER	2.4
1	A	92	ILE	2.4
1	B	209	ASP	2.4
1	A	95	MET	2.4
1	A	265	LEU	2.4
1	B	95	MET	2.4
1	B	149	GLY	2.3
1	C	99	SER	2.3
1	B	173	ASP	2.3
1	B	262	ALA	2.3
1	C	262	ALA	2.3
1	A	130	ILE	2.3
1	A	81	ILE	2.2
1	B	228	LEU	2.2
1	C	91	GLY	2.2
1	A	104	PRO	2.2
1	C	22	SER	2.2
1	A	269	ALA	2.2
1	B	259	HIS	2.1
1	B	260	SER	2.1
1	C	265	LEU	2.1
1	A	97	MET	2.1
1	B	110	ALA	2.1
1	C	175	GLY	2.1
1	B	94	LEU	2.1
1	A	79	VAL	2.0
1	A	172[A]	VAL	2.0
1	C	176	LEU	2.0
1	C	100	ASP	2.0
1	C	239	SER	2.0
1	B	67	PHE	2.0
1	B	70	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	CL	A	1290	1/1	0.91	0.12	52,52,52,52	0
2	CL	C	1284	1/1	0.92	0.12	57,57,57,57	0
2	CL	B	1283	1/1	0.93	0.31	30,30,30,30	0
2	CL	C	1283	1/1	0.96	0.09	49,49,49,49	0
2	CL	A	1291	1/1	0.99	0.51	30,30,30,30	0
2	CL	B	1284	1/1	0.99	0.51	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.