



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:24 AM UTC

PDB ID : 3ZVT / pdb_00003zvt
Title : Unexpected tricovalent binding mode of boronic acids within the active site of a penicillin binding protein
Authors : Sauvage, E.; Zervosen, A.; Herman, R.; Kerff, F.; Rocaboy, M.; Charlier, P.
Deposited on : 2011-07-27
Resolution : 3.10 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

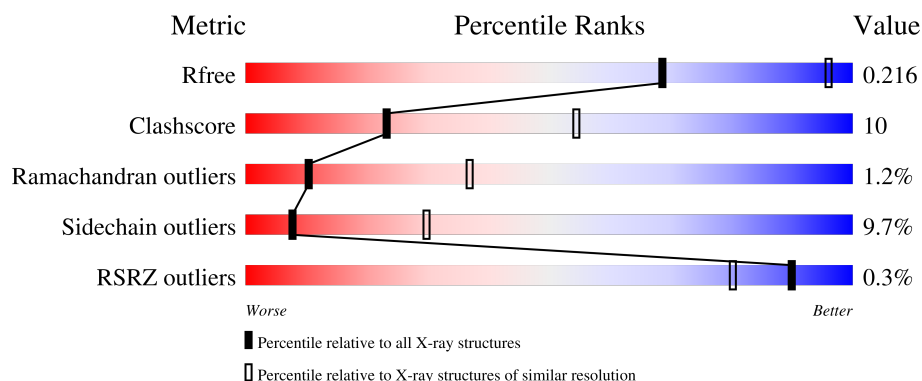
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	466	 74% 24% 2% 0%
1	B	466	 72% 26% 2% 0%
1	C	466	 70% 25% 5% 0%
1	D	466	 75% 22% 3% 0%

2 Entry composition [i](#)

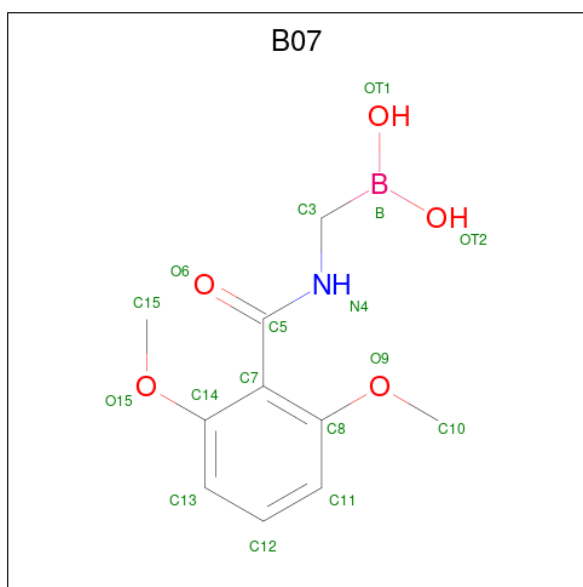
There are 4 unique types of molecules in this entry. The entry contains 13538 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 D-ALANYL-D-ALANINE CARBOXYPEPTIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	0	0	0
			3353	2076	564	707	6			
1	B	466	Total	C	N	O	S	0	0	1
			3344	2071	564	703	6			
1	C	466	Total	C	N	O	S	0	0	1
			3344	2071	564	703	6			
1	D	466	Total	C	N	O	S	0	0	0
			3353	2076	564	707	6			

- Molecule 2 is 2,6 DIMETHOXYBENZAMIDOBORONIC ACID (CCD ID: B07) (formula: $C_{10}H_{14}BNO_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	B	C	N	O	0
			15	1	10	1	3	0
2	B	1	Total	B	C	N	O	0
			15	1	10	1	3	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	B	C	N	O	0	0
			15	1	10	1	3		
2	D	1	Total	B	C	N	O	0	0
			15	1	10	1	3		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

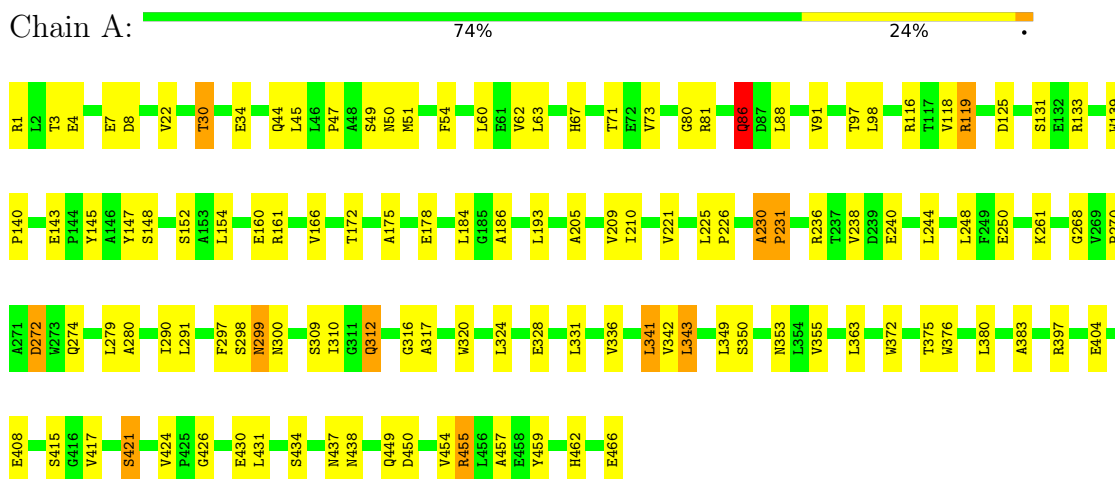
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		
4	D	2	Total	Mg	0	0
			2	2		

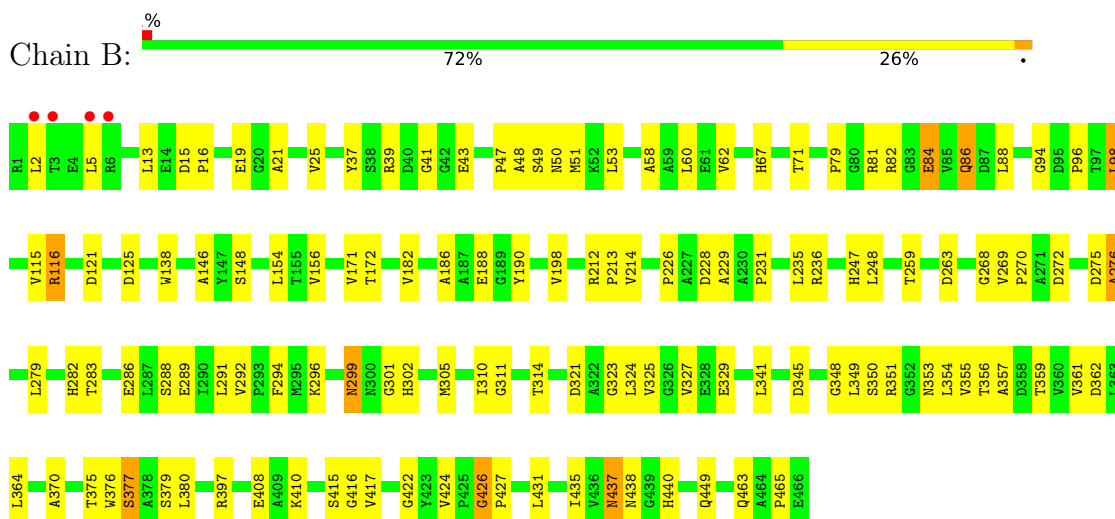
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE

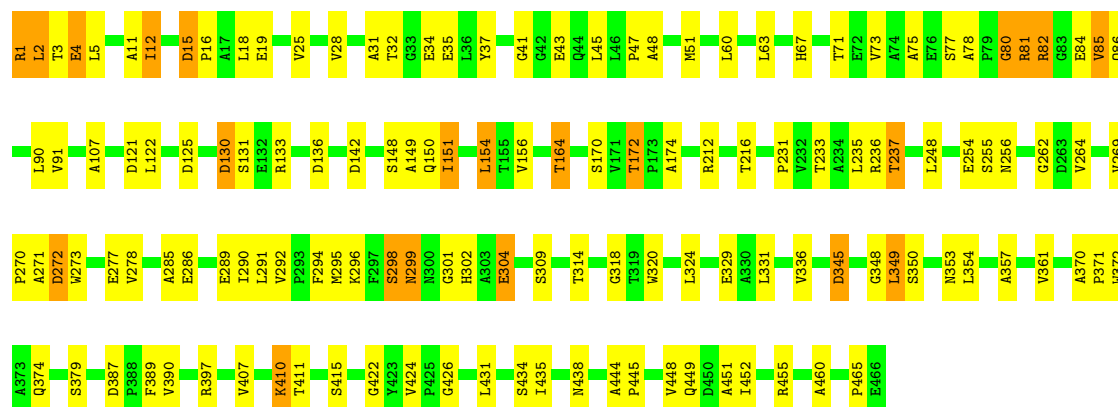


• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE



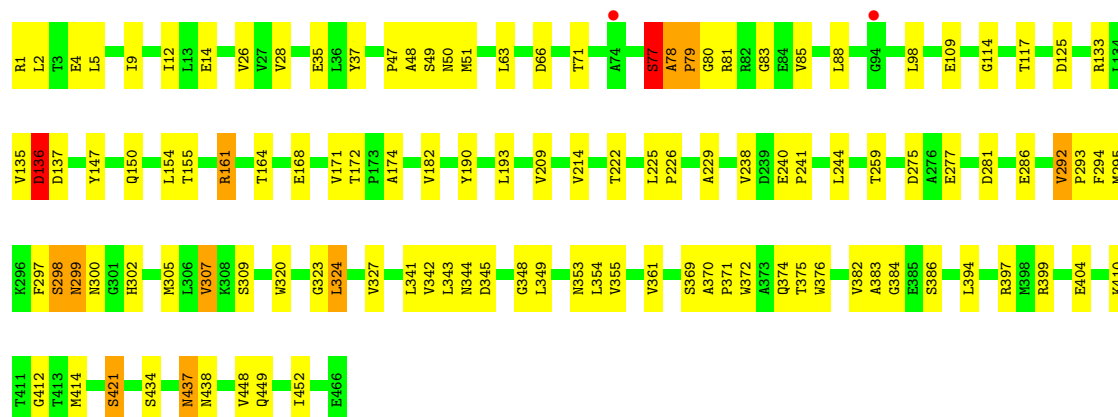
• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE





• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE

Chain D: 75% 22% •



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	103.99Å 92.23Å 107.04Å 90.00° 94.43° 90.00°	Depositor
Resolution (Å)	35.83 – 3.10 35.83 – 3.10	Depositor EDS
% Data completeness (in resolution range)	95.5 (35.83-3.10) 95.4 (35.83-3.10)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.203 , 0.254 0.203 , 0.216	Depositor DCC
R_{free} test set	1851 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	59.9	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13538	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.15% 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: SO4, B07, MG

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	0.60	0/3412	0.93	8/4666 (0.2%)
1	B	0.56	1/3403 (0.0%)	0.92	6/4656 (0.1%)
1	C	0.59	1/3403 (0.0%)	0.95	7/4656 (0.2%)
1	D	0.58	0/3412	0.91	4/4666 (0.1%)
All	All	0.58	2/13630 (0.0%)	0.93	25/18644 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	465	PRO	C-N	-6.64	1.24	1.33
1	C	465	PRO	C-N	-6.55	1.24	1.33

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	78	ALA	CA-C-N	6.82	128.37	119.84
1	D	78	ALA	C-N-CA	6.82	128.37	119.84
1	C	31	ALA	N-CA-C	-6.61	107.08	114.62
1	A	116	ARG	N-CA-C	-6.41	106.42	114.56
1	B	269	VAL	CA-C-N	6.04	127.39	119.84
1	B	269	VAL	C-N-CA	6.04	127.39	119.84
1	B	426	GLY	CA-C-N	5.82	125.49	119.56
1	B	426	GLY	C-N-CA	5.82	125.49	119.56
1	A	230	ALA	CA-C-N	5.81	127.10	119.84
1	A	230	ALA	C-N-CA	5.81	127.10	119.84
1	C	71	THR	N-CA-C	-5.72	100.36	109.23
1	C	15	ASP	CA-C-N	5.64	126.89	119.84
1	C	15	ASP	C-N-CA	5.64	126.89	119.84
1	A	172	THR	CA-C-N	-5.36	115.06	120.31
1	A	172	THR	C-N-CA	-5.36	115.06	120.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	166	VAL	N-CA-C	5.34	116.54	108.96
1	C	82	ARG	N-CA-C	-5.29	106.61	113.16
1	B	272	ASP	N-CA-C	-5.21	106.92	113.28
1	B	345	ASP	N-CA-C	5.18	115.89	108.74
1	D	109	GLU	N-CA-C	-5.18	107.09	113.41
1	D	137	ASP	N-CA-C	5.17	118.42	112.57
1	C	422	GLY	N-CA-C	5.08	117.70	110.74
1	A	426	GLY	CA-C-N	5.04	125.31	119.47
1	A	426	GLY	C-N-CA	5.04	125.31	119.47
1	C	81	ARG	N-CA-C	5.02	121.48	110.80

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	3353	0	3199	62	0
1	B	3344	0	3193	73	0
1	C	3344	0	3193	77	0
1	D	3353	0	3199	61	0
2	A	15	0	12	3	0
2	B	15	0	12	1	0
2	C	15	0	12	3	0
2	D	15	0	12	4	0
3	A	20	0	0	0	0
3	B	20	0	0	0	0
3	C	20	0	0	0	0
3	D	20	0	0	1	0
4	A	2	0	0	0	0
4	D	2	0	0	0	0
All	All	13538	0	12832	272	0

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 10.

All (272) 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:B:299:ASN:ND2	1:B:302:HIS:H	1.44	1.15
1:A:125:ASP:OD1	1:A:268:GLY:HA2	1.52	1.09
1:D:49:SER:HB3	1:D:412:GLY:HA2	1.49	0.94
1:B:60:LEU:HD21	1:B:291:LEU:HD11	1.53	0.91
1:C:149:ALA:HA	1:C:237:THR:HG21	1.57	0.87
1:A:51:MET:HE3	1:A:353:ASN:HB3	1.59	0.84
1:B:299:ASN:ND2	1:B:302:HIS:N	2.28	0.81
1:D:298:SER:CB	2:D:500:B07:C3	2.58	0.80
1:C:397:ARG:HH22	1:C:449:GLN:HE21	1.26	0.80
1:D:226:PRO:HG2	1:D:229:ALA:HB2	1.62	0.80
1:C:397:ARG:NH2	1:C:449:GLN:HE21	1.79	0.79
1:D:298:SER:CB	2:D:500:B07:H31C	2.14	0.78
1:C:397:ARG:HH22	1:C:449:GLN:NE2	1.83	0.77
1:B:292:VAL:O	1:B:296:LYS:HB2	1.84	0.77
1:A:50:ASN:ND2	1:A:421:SER:OG	2.18	0.76
1:C:353:ASN:O	1:C:354:LEU:HD23	1.86	0.76
1:C:2:LEU:C	1:C:4:GLU:H	1.93	0.76
1:A:457:ALA:O	1:A:462:HIS:HB2	1.86	0.76
1:B:288:SER:HB2	1:B:375:THR:HG21	1.67	0.75
1:D:49:SER:HB3	1:D:412:GLY:CA	2.17	0.74
1:B:408:GLU:O	1:B:422:GLY:HA3	1.86	0.74
1:B:299:ASN:HD22	1:B:302:HIS:H	1.33	0.73
1:A:91:VAL:HG22	1:A:125:ASP:HB3	1.69	0.73
1:C:270:PRO:C	1:C:272:ASP:H	1.97	0.72
1:A:97:THR:HG21	1:A:290:ILE:HG12	1.71	0.71
1:C:156:VAL:HG21	1:C:248:LEU:HD12	1.72	0.71
1:C:75:ALA:HB1	1:C:85:VAL:HG11	1.73	0.70
1:D:26:VAL:HG23	1:D:361:VAL:HG21	1.74	0.70
1:B:171:VAL:HG22	1:B:182:VAL:HG22	1.74	0.69
1:C:18:LEU:HD11	1:C:448:VAL:HG11	1.74	0.69
1:B:51:MET:HE3	1:B:353:ASN:CG	2.18	0.69
1:B:125:ASP:OD2	1:B:268:GLY:HA2	1.92	0.69
1:D:342:VAL:HB	1:D:354:LEU:HB2	1.74	0.68
1:A:147:TYR:HB2	1:A:300:ASN:ND2	2.09	0.68
1:C:47:PRO:HG3	1:C:51:MET:CE	2.24	0.68
1:B:341:LEU:HD23	1:B:355:VAL:HG12	1.76	0.68
1:D:79:PRO:HB3	1:D:114:GLY:HA3	1.75	0.67
1:D:50:ASN:ND2	1:D:421:SER:OG	2.27	0.67
1:D:421:SER:HB3	1:D:434:SER:HA	1.74	0.67
1:C:296:LYS:HE2	1:C:389:PHE:O	1.95	0.66
1:B:361:VAL:HA	1:B:364:LEU:HD12	1.76	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:GLY:HA2	1:B:314:THR:O	1.96	0.65
1:A:238:VAL:HG21	1:A:244:LEU:HD22	1.79	0.65
1:C:361:VAL:HG23	1:C:434:SER:HB3	1.80	0.64
1:B:377:SER:HA	1:B:380:LEU:HD12	1.79	0.63
1:C:349:LEU:HD23	2:C:500:B07:H13	1.80	0.63
1:C:48:ALA:O	1:C:348:GLY:HA3	1.99	0.62
1:A:73:VAL:HG13	1:A:88:LEU:HD21	1.80	0.62
1:C:80:GLY:O	1:C:82:ARG:N	2.22	0.62
1:C:410:LYS:HG3	1:C:411:THR:N	2.13	0.62
1:B:172:THR:HG22	1:B:231:PRO:HB3	1.82	0.62
1:D:298:SER:OG	2:D:500:B07:H31C	1.82	0.62
1:A:51:MET:CE	1:A:353:ASN:HB3	2.30	0.62
1:C:294:PHE:HB2	1:C:302:HIS:HD2	1.64	0.62
1:A:47:PRO:HG3	1:A:51:MET:HE2	1.81	0.61
1:B:351:ARG:NH2	1:B:415:SER:O	2.33	0.61
1:A:175:ALA:HB3	1:A:178:GLU:OE2	2.01	0.61
1:C:331:LEU:O	1:C:336:VAL:HG12	2.01	0.61
1:B:86:GLN:HA	1:B:86:GLN:OE1	2.01	0.60
1:A:147:TYR:HB2	1:A:300:ASN:HD22	1.65	0.60
1:A:298:SER:CB	2:A:500:B07:C3	2.79	0.60
1:D:448:VAL:O	1:D:452:ILE:HG13	2.00	0.60
1:B:294:PHE:HB2	1:B:302:HIS:HD2	1.67	0.60
1:A:397:ARG:HH22	1:A:449:GLN:HE21	1.50	0.59
1:B:172:THR:CG2	1:B:231:PRO:HB3	2.32	0.59
1:C:25:VAL:HG22	1:C:435:ILE:HG23	1.84	0.59
1:C:172:THR:HG23	1:C:231:PRO:HB3	1.84	0.59
1:A:119:ARG:HE	1:A:261:LYS:HE2	1.67	0.59
1:A:421:SER:HB3	1:A:434:SER:HA	1.84	0.58
1:D:294:PHE:HB2	1:D:302:HIS:HD2	1.67	0.58
1:B:49:SER:CB	2:B:500:B07:C3	2.81	0.58
1:D:383:ALA:HA	1:D:394:LEU:HB3	1.84	0.58
1:C:16:PRO:O	1:C:19:GLU:HB2	2.04	0.58
1:A:331:LEU:O	1:A:336:VAL:HG12	2.04	0.58
1:C:5:LEU:HA	1:C:455:ARG:HH12	1.68	0.58
1:C:164:THR:HG21	1:C:301:GLY:HA3	1.85	0.57
1:C:28:VAL:HG12	1:C:35:GLU:HA	1.85	0.57
1:D:150:GLN:NE2	1:D:240:GLU:H	2.02	0.57
1:C:130:ASP:HB3	1:C:318:GLY:O	2.03	0.57
1:D:155:THR:OG1	1:D:241:PRO:HB2	2.04	0.57
1:B:82:ARG:HB2	1:B:84:GLU:OE1	2.05	0.56
1:B:51:MET:CE	1:B:353:ASN:HB3	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:51:MET:HE3	1:D:353:ASN:CG	2.30	0.56
1:D:77:SER:HA	1:D:277:GLU:HG3	1.87	0.56
1:D:1:ARG:HA	1:D:4:GLU:HB2	1.87	0.56
1:B:299:ASN:HD21	1:B:302:HIS:H	1.44	0.56
1:D:370:ALA:O	1:D:372:TRP:N	2.38	0.56
1:B:16:PRO:HA	1:B:19:GLU:HG2	1.86	0.56
1:B:437:ASN:C	1:B:438:ASN:HD22	2.14	0.56
1:B:79:PRO:HG3	1:B:115:VAL:HG22	1.87	0.55
1:C:151:ILE:HG12	1:C:304:GLU:HB3	1.89	0.55
1:B:299:ASN:HD22	1:B:299:ASN:C	2.14	0.55
1:D:49:SER:CB	1:D:412:GLY:HA2	2.32	0.55
1:D:150:GLN:HE22	1:D:240:GLU:H	1.54	0.55
1:B:212:ARG:O	1:B:213:PRO:C	2.50	0.55
1:B:235:LEU:O	1:B:236:ARG:HD3	2.07	0.55
1:A:60:LEU:HD13	1:A:372:TRP:HB2	1.89	0.55
1:C:426:GLY:HA3	1:C:460:ALA:HB1	1.88	0.55
1:D:348:GLY:HA2	1:D:353:ASN:HD22	1.71	0.55
1:C:149:ALA:HA	1:C:237:THR:CG2	2.35	0.54
1:D:135:VAL:O	1:D:136:ASP:C	2.51	0.54
1:B:156:VAL:HG21	1:B:248:LEU:HD12	1.90	0.53
1:C:121:ASP:HA	1:C:262:GLY:HA3	1.90	0.53
1:D:225:LEU:HD12	1:D:226:PRO:HD2	1.90	0.53
1:C:122:LEU:HD23	1:C:264:VAL:HG22	1.90	0.53
1:A:131:SER:HB2	1:A:133:ARG:HH21	1.74	0.53
1:B:43:GLU:O	1:B:356:THR:HA	2.08	0.53
1:B:71:THR:HG21	1:B:98:LEU:HD11	1.90	0.53
1:D:28:VAL:HG23	1:D:35:GLU:HA	1.91	0.53
1:C:151:ILE:CG1	1:C:304:GLU:HB3	2.39	0.53
1:A:298:SER:CB	2:A:500:B07:H31C	2.39	0.52
1:B:37:TYR:CE2	1:B:39:ARG:HB3	2.45	0.52
1:B:294:PHE:HB2	1:B:302:HIS:CD2	2.45	0.52
1:B:416:GLY:HA2	1:B:440:HIS:CD2	2.45	0.52
1:C:298:SER:CB	2:C:500:B07:C3	2.88	0.52
1:B:397:ARG:HH22	1:B:449:GLN:NE2	2.08	0.52
1:D:299:ASN:HB3	1:D:302:HIS:HB2	1.91	0.52
1:A:30:THR:HB	1:A:430:GLU:HB3	1.93	0.51
1:C:292:VAL:O	1:C:296:LYS:HG2	2.11	0.51
1:D:117:THR:HG23	1:D:259:THR:HB	1.93	0.51
1:D:190:TYR:CE2	1:D:244:LEU:HB2	2.45	0.51
1:A:184:LEU:HD21	1:A:193:LEU:HD13	1.92	0.51
1:D:26:VAL:HG23	1:D:361:VAL:CG2	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:142:ASP:O	1:C:148:SER:HB3	2.10	0.51
1:D:345:ASP:C	1:D:345:ASP:OD1	2.54	0.51
1:A:91:VAL:HG22	1:A:125:ASP:CB	2.38	0.51
1:A:270:PRO:C	1:A:272:ASP:H	2.19	0.51
1:B:47:PRO:HG3	1:B:51:MET:HE2	1.93	0.51
1:B:275:ASP:O	1:B:276:ALA:C	2.54	0.51
1:B:41:GLY:HA2	1:B:357:ALA:HB3	1.94	0.50
1:C:63:LEU:O	1:C:67:HIS:HB2	2.12	0.50
1:A:312:GLN:O	1:A:316:GLY:N	2.42	0.50
1:C:345:ASP:OD1	1:C:350:SER:HB3	2.12	0.49
1:B:25:VAL:HG22	1:B:435:ILE:HG23	1.94	0.49
1:A:139:TRP:N	1:A:139:TRP:CD1	2.79	0.49
1:B:79:PRO:HD3	1:B:279:LEU:HD21	1.95	0.49
1:C:164:THR:CG2	1:C:301:GLY:HA3	2.43	0.49
1:C:424:VAL:HB	1:C:431:LEU:HB2	1.94	0.49
1:D:370:ALA:C	1:D:372:TRP:H	2.20	0.49
1:C:285:ALA:O	1:C:290:ILE:HD11	2.13	0.49
1:B:67:HIS:O	1:B:286:GLU:HA	2.13	0.49
1:C:41:GLY:HA2	1:C:357:ALA:HB3	1.94	0.48
1:C:47:PRO:HG3	1:C:51:MET:HE2	1.95	0.48
1:D:47:PRO:O	1:D:50:ASN:HB2	2.13	0.48
1:A:49:SER:N	2:A:500:B07:H151	2.27	0.48
1:B:138:TRP:HE1	1:B:350:SER:HB2	1.79	0.48
1:C:448:VAL:O	1:C:451:ALA:HB3	2.13	0.48
1:D:307:VAL:HG13	1:D:323:GLY:HA3	1.96	0.48
1:B:94:GLY:O	1:B:96:PRO:HD3	2.13	0.48
1:B:96:PRO:HB2	1:B:302:HIS:CE1	2.49	0.48
1:D:48:ALA:O	1:D:348:GLY:HA3	2.13	0.48
1:D:133:ARG:N	3:D:603:SO4:O3	2.45	0.48
1:A:350:SER:O	1:A:353:ASN:ND2	2.44	0.48
1:A:54:PHE:CD1	1:A:363:LEU:HD22	2.49	0.48
1:C:397:ARG:NH2	1:C:449:GLN:NE2	2.49	0.48
1:B:116:ARG:HB3	1:B:116:ARG:NH1	2.29	0.47
1:A:417:VAL:HG13	1:A:438:ASN:HD21	1.78	0.47
1:C:90:LEU:HD23	1:C:154:LEU:HD11	1.95	0.47
1:C:448:VAL:O	1:C:452:ILE:HD12	2.13	0.47
1:C:407:VAL:HG22	1:C:424:VAL:HG22	1.97	0.47
1:C:298:SER:CB	2:C:500:B07:H31C	2.44	0.47
1:A:383:ALA:O	1:A:404:GLU:HA	2.14	0.47
1:B:397:ARG:HH22	1:B:449:GLN:HE21	1.63	0.47
1:D:171:VAL:HG22	1:D:182:VAL:HG22	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:172:THR:CG2	1:C:231:PRO:HB3	2.44	0.47
1:C:299:ASN:HB3	1:C:302:HIS:HB2	1.97	0.47
1:B:146:ALA:C	1:B:148:SER:H	2.23	0.47
1:B:301:GLY:O	1:B:305:MET:HG3	2.14	0.47
1:A:145:TYR:HB2	1:A:147:TYR:CE1	2.49	0.46
1:B:424:VAL:HB	1:B:431:LEU:HB2	1.97	0.46
1:D:14:GLU:OE1	1:D:14:GLU:HA	2.16	0.46
1:D:384:GLY:HA3	1:D:404:GLU:HG3	1.97	0.46
1:A:230:ALA:HA	1:A:231:PRO:HD2	1.75	0.46
1:B:310:ILE:HD13	1:B:327:VAL:HG22	1.97	0.46
1:C:15:ASP:HA	1:C:16:PRO:HD2	1.75	0.46
1:C:270:PRO:C	1:C:272:ASP:N	2.64	0.45
1:A:47:PRO:O	1:A:50:ASN:HB2	2.16	0.45
1:A:424:VAL:O	1:A:431:LEU:HB2	2.16	0.45
1:B:226:PRO:HG2	1:B:229:ALA:HB2	1.97	0.45
1:A:73:VAL:CG1	1:A:88:LEU:HD21	2.47	0.45
1:D:81:ARG:C	1:D:83:GLY:H	2.25	0.45
1:D:47:PRO:HG3	1:D:51:MET:HE2	1.99	0.45
1:D:161:ARG:HG2	1:D:297:PHE:CD2	2.52	0.45
1:A:51:MET:HE3	1:A:353:ASN:CB	2.39	0.45
1:C:91:VAL:HG22	1:C:125:ASP:HB3	1.98	0.45
1:C:444:ALA:HA	1:C:445:PRO:HD3	1.80	0.45
1:D:397:ARG:HH22	1:D:449:GLN:HE21	1.64	0.45
1:A:62:VAL:HG11	1:A:310:ILE:HG23	1.99	0.44
1:C:387:ASP:OD2	1:C:390:VAL:HG23	2.18	0.44
1:D:147:TYR:HB2	1:D:300:ASN:ND2	2.33	0.44
1:A:1:ARG:NH2	1:A:459:TYR:HA	2.32	0.44
1:A:299:ASN:C	1:A:299:ASN:HD22	2.25	0.44
1:A:63:LEU:O	1:A:67:HIS:HB2	2.17	0.44
1:D:437:ASN:C	1:D:438:ASN:HD22	2.25	0.44
1:A:86:GLN:O	1:A:118:VAL:HG13	2.18	0.44
1:A:417:VAL:HG13	1:A:438:ASN:ND2	2.32	0.44
1:B:21:ALA:HB2	1:B:440:HIS:HA	2.00	0.44
1:B:96:PRO:HB2	1:B:302:HIS:ND1	2.33	0.44
1:A:45:LEU:HB3	1:A:438:ASN:OD1	2.18	0.44
1:A:186:ALA:HB1	1:A:248:LEU:HD11	1.99	0.44
1:B:190:TYR:HB2	1:B:247:HIS:CD2	2.53	0.44
1:C:212:ARG:HD2	1:C:216:THR:O	2.18	0.44
1:C:370:ALA:HB1	1:C:371:PRO:HD2	1.99	0.44
1:A:125:ASP:OD1	1:A:268:GLY:CA	2.44	0.43
1:B:49:SER:OG	1:B:410:LYS:CE	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:270:PRO:HG2	1:C:273:TRP:CE2	2.53	0.43
1:A:161:ARG:HG2	1:A:297:PHE:CE2	2.52	0.43
1:B:299:ASN:ND2	1:B:299:ASN:C	2.76	0.43
1:C:235:LEU:O	1:C:236:ARG:NH1	2.49	0.43
1:D:323:GLY:O	1:D:327:VAL:HG23	2.18	0.43
1:D:294:PHE:CD1	1:D:302:HIS:HB3	2.54	0.43
1:A:8:ASP:OD2	1:A:455:ARG:NH1	2.49	0.43
1:C:1:ARG:HG2	1:C:2:LEU:HG	2.01	0.43
1:A:140:PRO:HA	1:A:143:GLU:HG3	2.00	0.43
1:C:73:VAL:HG22	1:C:90:LEU:HD12	2.00	0.43
1:C:133:ARG:HB3	1:C:150:GLN:HG2	2.01	0.43
1:C:78:ALA:N	1:C:277:GLU:HG2	2.33	0.43
1:C:170:SER:HB3	1:C:233:THR:HG23	2.00	0.43
1:D:292:VAL:HB	1:D:293:PRO:CD	2.49	0.43
1:B:53:LEU:O	1:B:376:TRP:HH2	2.02	0.43
1:B:296:LYS:HD2	1:B:379:SER:O	2.19	0.43
1:B:323:GLY:O	1:B:327:VAL:HG23	2.19	0.43
1:C:142:ASP:O	1:C:148:SER:CB	2.66	0.42
1:C:345:ASP:OD1	1:C:353:ASN:ND2	2.52	0.42
1:B:13:LEU:C	1:B:15:ASP:H	2.26	0.42
1:D:298:SER:HB2	2:D:500:B07:H31C	1.96	0.42
1:A:143:GLU:HA	1:A:148:SER:OG	2.19	0.42
1:C:45:LEU:HD22	1:C:438:ASN:HB2	1.99	0.42
1:C:372:TRP:HE3	1:C:372:TRP:O	2.02	0.42
1:D:66:ASP:HA	1:D:286:GLU:OE1	2.18	0.42
1:B:228:ASP:OD1	1:B:228:ASP:N	2.52	0.42
1:B:361:VAL:O	1:B:362:ASP:C	2.63	0.42
1:A:73:VAL:HB	1:A:280:ALA:HB3	2.02	0.42
1:A:341:LEU:HD22	1:A:343:LEU:HD22	2.00	0.42
1:D:294:PHE:CD2	1:D:295:MET:HG2	2.54	0.42
1:D:320:TRP:O	1:D:324:LEU:HD22	2.18	0.42
1:B:48:ALA:HA	1:B:348:GLY:O	2.20	0.42
1:A:210:ILE:HB	1:A:236:ARG:NH2	2.35	0.42
1:B:341:LEU:CD2	1:B:355:VAL:HG12	2.47	0.42
1:C:107:ALA:CB	1:C:256:ASN:HD22	2.32	0.42
1:B:282:HIS:CE1	1:B:283:THR:O	2.72	0.42
1:C:331:LEU:C	1:C:336:VAL:HG12	2.44	0.42
1:D:369:SER:O	1:D:370:ALA:C	2.63	0.42
1:D:12:ILE:HG21	1:D:452:ILE:HG12	2.02	0.42
1:D:341:LEU:HD13	1:D:343:LEU:HD11	2.01	0.42
1:B:311:GLY:CA	1:B:323:GLY:HA2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:LEU:HG	1:B:359:THR:HG21	2.01	0.41
1:C:51:MET:HE3	1:C:353:ASN:CG	2.45	0.41
1:C:60:LEU:HD13	1:C:372:TRP:CE3	2.55	0.41
1:C:286:GLU:HB2	1:C:289:GLU:HG2	2.03	0.41
1:A:225:LEU:HA	1:A:226:PRO:HD3	1.91	0.41
1:C:11:ALA:O	1:C:12:ILE:C	2.64	0.41
1:D:77:SER:HB2	1:D:78:ALA:H	1.66	0.41
1:D:26:VAL:HA	1:D:37:TYR:O	2.21	0.41
1:D:79:PRO:HB2	1:D:80:GLY:H	1.69	0.41
1:A:47:PRO:CG	1:A:51:MET:HE2	2.50	0.40
1:B:186:ALA:C	1:B:188:GLU:H	2.28	0.40
1:A:240:GLU:OE1	1:A:240:GLU:HA	2.21	0.40
1:C:295:MET:HB2	1:C:379:SER:HB2	2.03	0.40
1:D:125:ASP:C	1:D:125:ASP:OD1	2.63	0.40
1:D:386:SER:HA	1:D:399:ARG:HH21	1.85	0.40
1:A:380:LEU:HB2	1:A:408:GLU:HB3	2.03	0.40
1:A:450:ASP:O	1:A:454:VAL:HG23	2.21	0.40
1:C:32:THR:C	1:C:34:GLU:N	2.79	0.40
1:B:426:GLY:HA2	1:B:427:PRO:HD3	1.90	0.40
1:C:130:ASP:OD1	1:C:131:SER:N	2.54	0.40
1:D:5:LEU:O	1:D:9:ILE:HG13	2.20	0.40
1:A:133:ARG:NH1	1:A:152:SER:HB2	2.37	0.40
1:A:312:GLN:HA	1:A:317:ALA:H	1.87	0.40
1:B:58:ALA:O	1:B:62:VAL:HG23	2.22	0.40
1:B:321:ASP:O	1:B:325:VAL:HG23	2.22	0.40
1:B:416:GLY:O	1:B:417:VAL:HG23	2.21	0.40

There are no symmetry-related clashes.

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	464/466 (100%)	418 (90%)	43 (9%)	3 (1%)	21	52
1	B	464/466 (100%)	411 (89%)	48 (10%)	5 (1%)	11	39
1	C	464/466 (100%)	422 (91%)	34 (7%)	8 (2%)	7	30
1	D	464/466 (100%)	416 (90%)	42 (9%)	6 (1%)	9	35
All	All	1856/1864 (100%)	1667 (90%)	167 (9%)	22 (1%)	10	37

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	214	VAL
1	B	370	ALA
1	C	81	ARG
1	D	79	PRO
1	B	276	ALA
1	C	80	GLY
1	C	37	TYR
1	D	371	PRO
1	A	86	GLN
1	A	205	ALA
1	B	86	GLN
1	C	86	GLN
1	C	174	ALA
1	D	136	ASP
1	D	174	ALA
1	A	231	PRO
1	C	3	THR
1	D	77	SER
1	D	292	VAL
1	C	271	ALA
1	C	12	ILE
1	B	270	PRO

5.3.2 Protein sidechains ⓘ

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	339/339 (100%)	300 (88%)	39 (12%)	5	23
1	B	338/339 (100%)	316 (94%)	22 (6%)	15	44
1	C	338/339 (100%)	305 (90%)	33 (10%)	7	29
1	D	339/339 (100%)	302 (89%)	37 (11%)	6	24
All	All	1354/1356 (100%)	1223 (90%)	131 (10%)	8	30

All (131) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	4	GLU
1	A	7	GLU
1	A	22	VAL
1	A	30	THR
1	A	34	GLU
1	A	44	GLN
1	A	71	THR
1	A	81	ARG
1	A	86	GLN
1	A	98	LEU
1	A	119	ARG
1	A	154	LEU
1	A	160	GLU
1	A	209	VAL
1	A	221	VAL
1	A	250	GLU
1	A	272	ASP
1	A	274	GLN
1	A	279	LEU
1	A	291	LEU
1	A	299	ASN
1	A	309	SER
1	A	312	GLN
1	A	320	TRP
1	A	324	LEU
1	A	328	GLU
1	A	341	LEU
1	A	342	VAL
1	A	343	LEU
1	A	349	LEU
1	A	355	VAL
1	A	375	THR

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Mol	Chain	Res	Type
1	A	376	TRP
1	A	415	SER
1	A	421	SER
1	A	437	ASN
1	A	455	ARG
1	A	466	GLU
1	B	2	LEU
1	B	5	LEU
1	B	50	ASN
1	B	81	ARG
1	B	84	GLU
1	B	88	LEU
1	B	98	LEU
1	B	116	ARG
1	B	121	ASP
1	B	154	LEU
1	B	198	VAL
1	B	259	THR
1	B	263	ASP
1	B	289	GLU
1	B	299	ASN
1	B	324	LEU
1	B	329	GLU
1	B	349	LEU
1	B	354	LEU
1	B	377	SER
1	B	437	ASN
1	B	463	GLN
1	C	1	ARG
1	C	2	LEU
1	C	4	GLU
1	C	43	GLU
1	C	77	SER
1	C	84	GLU
1	C	85	VAL
1	C	130	ASP
1	C	136	ASP
1	C	151	ILE
1	C	154	LEU
1	C	164	THR
1	C	172	THR
1	C	237	THR

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Mol	Chain	Res	Type
1	C	254	GLU
1	C	255	SER
1	C	269	VAL
1	C	272	ASP
1	C	278	VAL
1	C	291	LEU
1	C	298	SER
1	C	299	ASN
1	C	304	GLU
1	C	309	SER
1	C	314	THR
1	C	320	TRP
1	C	324	LEU
1	C	329	GLU
1	C	345	ASP
1	C	349	LEU
1	C	374	GLN
1	C	410	LYS
1	C	415	SER
1	D	2	LEU
1	D	63	LEU
1	D	71	THR
1	D	77	SER
1	D	85	VAL
1	D	88	LEU
1	D	98	LEU
1	D	136	ASP
1	D	154	LEU
1	D	161	ARG
1	D	164	THR
1	D	168	GLU
1	D	172	THR
1	D	193	LEU
1	D	209	VAL
1	D	214	VAL
1	D	222	THR
1	D	238	VAL
1	D	275	ASP
1	D	281	ASP
1	D	298	SER
1	D	299	ASN
1	D	305	MET

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Mol	Chain	Res	Type
1	D	307	VAL
1	D	309	SER
1	D	324	LEU
1	D	344	ASN
1	D	349	LEU
1	D	355	VAL
1	D	374	GLN
1	D	375	THR
1	D	376	TRP
1	D	382	VAL
1	D	410	LYS
1	D	414	MET
1	D	421	SER
1	D	437	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	50	ASN
1	A	274	GLN
1	A	282	HIS
1	A	299	ASN
1	A	312	GLN
1	A	449	GLN
1	A	463	GLN
1	B	50	ASN
1	B	67	HIS
1	B	150	GLN
1	B	158	HIS
1	B	299	ASN
1	B	302	HIS
1	B	396	ASN
1	B	437	ASN
1	B	449	GLN
1	C	50	ASN
1	C	247	HIS
1	C	256	ASN
1	C	299	ASN
1	C	302	HIS
1	C	449	GLN
1	D	50	ASN

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Mol	Chain	Res	Type
1	D	67	HIS
1	D	150	GLN
1	D	158	HIS
1	D	299	ASN
1	D	302	HIS
1	D	344	ASN
1	D	449	GLN
1	D	462	HIS

5.3.3 RNA [i](#)

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 24 ligands modelled in this entry, 4 are monoatomic - leaving 20 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	D	601	-	4,4,4	0.31	0	6,6,6	0.18	0
3	SO4	D	602	-	4,4,4	0.24	0	6,6,6	0.20	0
3	SO4	C	604	-	4,4,4	0.23	0	6,6,6	0.12	0
3	SO4	B	603	-	4,4,4	0.24	0	6,6,6	0.22	0
3	SO4	B	601	-	4,4,4	0.26	0	6,6,6	0.18	0
3	SO4	B	602	-	4,4,4	0.25	0	6,6,6	0.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	D	603	-	4,4,4	0.26	0	6,6,6	0.24	0
3	SO4	A	603	-	4,4,4	0.22	0	6,6,6	0.14	0
3	SO4	D	604	-	4,4,4	0.24	0	6,6,6	0.11	0
3	SO4	A	604	-	4,4,4	0.26	0	6,6,6	0.30	0
2	B07	C	500	1	14,15,17	0.57	0	18,19,22	2.08	7 (38%)
3	SO4	A	601	-	4,4,4	0.23	0	6,6,6	0.22	0
3	SO4	B	604	-	4,4,4	0.21	0	6,6,6	0.16	0
3	SO4	C	603	-	4,4,4	0.28	0	6,6,6	0.25	0
3	SO4	A	602	-	4,4,4	0.23	0	6,6,6	0.13	0
2	B07	D	500	1	14,15,17	0.60	0	18,19,22	1.66	6 (33%)
3	SO4	C	601	-	4,4,4	0.30	0	6,6,6	0.18	0
2	B07	A	500	1	14,15,17	0.46	0	18,19,22	1.99	6 (33%)
2	B07	B	500	1	14,15,17	0.51	0	18,19,22	1.85	8 (44%)
3	SO4	C	602	-	4,4,4	0.27	0	6,6,6	0.12	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	B07	A	500	1	-	2/10/11/13	0/1/1/1
2	B07	B	500	1	-	2/10/11/13	0/1/1/1
2	B07	C	500	1	-	4/10/11/13	0/1/1/1
2	B07	D	500	1	-	4/10/11/13	0/1/1/1

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	B07	C10-O9-C8	4.82	124.58	117.51
2	A	500	B07	O9-C8-C7	4.03	121.55	115.84
2	C	500	B07	C15-O15-C14	3.79	123.07	117.51
2	A	500	B07	O9-C8-C11	-3.36	118.64	124.30
2	C	500	B07	O9-C8-C7	3.21	120.39	115.84
2	D	500	B07	O9-C8-C7	3.14	120.28	115.84
2	A	500	B07	C7-C5-N4	2.94	121.40	115.67
2	B	500	B07	C15-O15-C14	2.92	121.79	117.51
2	A	500	B07	C10-O9-C8	2.89	121.76	117.51
2	D	500	B07	O15-C14-C7	2.88	119.92	115.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	B07	O15-C14-C7	2.80	119.81	115.84
2	D	500	B07	O15-C14-C13	-2.73	119.70	124.30
2	B	500	B07	C10-O9-C8	2.70	121.47	117.51
2	B	500	B07	O15-C14-C13	-2.68	119.79	124.30
2	B	500	B07	O9-C8-C7	2.59	119.51	115.84
2	D	500	B07	O9-C8-C11	-2.42	120.23	124.30
2	B	500	B07	O9-C8-C11	-2.36	120.32	124.30
2	C	500	B07	C8-C7-C14	2.36	121.36	117.83
2	B	500	B07	C7-C5-N4	2.31	120.16	115.67
2	C	500	B07	O15-C14-C13	-2.28	120.46	124.30
2	C	500	B07	C7-C5-N4	2.25	120.05	115.67
2	C	500	B07	O15-C14-C7	2.21	118.98	115.84
2	A	500	B07	C15-O15-C14	2.19	120.73	117.51
2	B	500	B07	C3-N4-C5	2.13	125.69	122.06
2	D	500	B07	C10-O9-C8	2.13	120.64	117.51
2	D	500	B07	C7-C5-N4	2.12	119.79	115.67
2	A	500	B07	C14-C7-C5	-2.04	116.79	120.66

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	B07	C11-C8-O9-C10
2	D	500	B07	C11-C8-O9-C10
2	B	500	B07	C11-C8-O9-C10
2	D	500	B07	C7-C8-O9-C10
2	B	500	B07	C7-C8-O9-C10
2	D	500	B07	C13-C14-O15-C15
2	A	500	B07	C7-C8-O9-C10
2	C	500	B07	C13-C14-O15-C15
2	C	500	B07	C11-C8-O9-C10
2	D	500	B07	C7-C14-O15-C15
2	C	500	B07	C7-C14-O15-C15
2	C	500	B07	C7-C8-O9-C10

There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	603	SO4	1	0
2	C	500	B07	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	500	B07	4	0
2	A	500	B07	3	0
2	B	500	B07	1	0

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	466/466 (100%)	-0.43	0 100 100	35, 55, 84, 114	0
1	B	466/466 (100%)	-0.41	4 (0%) 81 63	33, 57, 90, 237	0
1	C	466/466 (100%)	-0.46	0 100 100	37, 55, 83, 121	0
1	D	466/466 (100%)	-0.20	2 (0%) 88 76	44, 66, 107, 137	0
All	All	1864/1864 (100%)	-0.38	6 (0%) 90 80	33, 58, 96, 237	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	5	LEU	2.9
1	B	3	THR	2.8
1	B	2	LEU	2.3
1	D	74	ALA	2.2
1	D	94	GLY	2.2
1	B	6	ARG	2.2

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

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(Å ²)	Q<0.9
3	SO4	D	604	5/5	0.67	0.13	151,151,151,151	0
3	SO4	A	604	5/5	0.70	0.14	117,117,117,117	0
3	SO4	A	602	5/5	0.80	0.12	112,112,112,113	0
3	SO4	B	602	5/5	0.81	0.13	98,98,98,99	0
3	SO4	B	604	5/5	0.82	0.12	114,114,114,115	0
3	SO4	D	601	5/5	0.83	0.11	91,91,92,92	0
3	SO4	C	604	5/5	0.83	0.13	116,116,116,116	0
3	SO4	C	602	5/5	0.84	0.10	98,99,99,99	0
3	SO4	A	603	5/5	0.88	0.19	119,119,120,120	0
3	SO4	D	602	5/5	0.89	0.10	86,86,86,87	0
3	SO4	C	603	5/5	0.91	0.09	83,83,84,84	0
3	SO4	D	603	5/5	0.92	0.14	75,75,76,76	0
3	SO4	A	601	5/5	0.92	0.08	64,64,64,64	0
4	MG	A	610	1/1	0.92	0.06	60,60,60,60	0
2	B07	D	500	15/17	0.93	0.09	40,48,50,51	0
3	SO4	B	601	5/5	0.93	0.10	77,77,77,78	0
2	B07	A	500	15/17	0.93	0.09	34,51,53,54	0
3	SO4	B	603	5/5	0.94	0.08	68,68,68,69	0
2	B07	C	500	15/17	0.95	0.09	30,42,44,45	0
3	SO4	C	601	5/5	0.95	0.07	63,64,64,65	0
4	MG	A	611	1/1	0.95	0.04	64,64,64,64	0
2	B07	B	500	15/17	0.96	0.07	31,43,46,47	0
4	MG	D	610	1/1	0.96	0.05	69,69,69,69	0
4	MG	D	611	1/1	0.99	0.04	35,35,35,35	0

6.5 Other polymers

There are no such residues in this entry.