



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 02:18 am GMT

PDB ID : 3ZVT
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

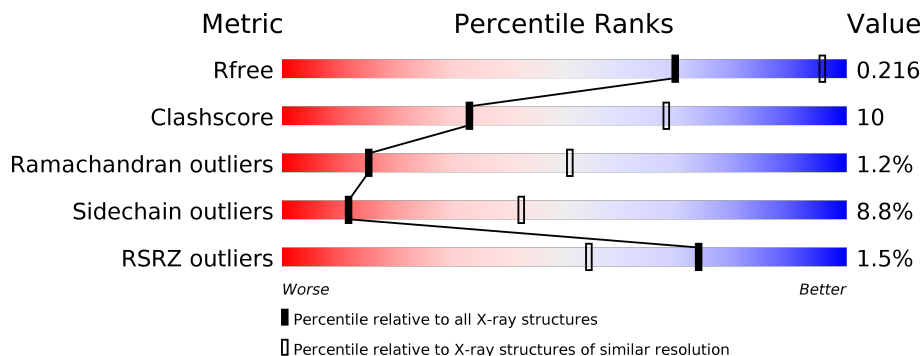
1 Overall quality at a glance

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}	100719	1001 (3.12-3.08)
Clashscore	112137	1099 (3.12-3.08)
Ramachandran outliers	110173	1057 (3.12-3.08)
Sidechain outliers	110143	1057 (3.12-3.08)
RSRZ outliers	101464	1006 (3.12-3.08)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>76%</div> <div>22%</div> <div>•</div> </div>
1	B	466	<div> <div>4%</div> <div>74%</div> <div>24%</div> <div>•</div> </div>
1	C	466	<div> <div>4%</div> <div>72%</div> <div>25%</div> <div>•</div> </div>
1	D	466	<div> <div>4%</div> <div>77%</div> <div>21%</div> <div>•</div> </div>

2 Entry composition [i](#)

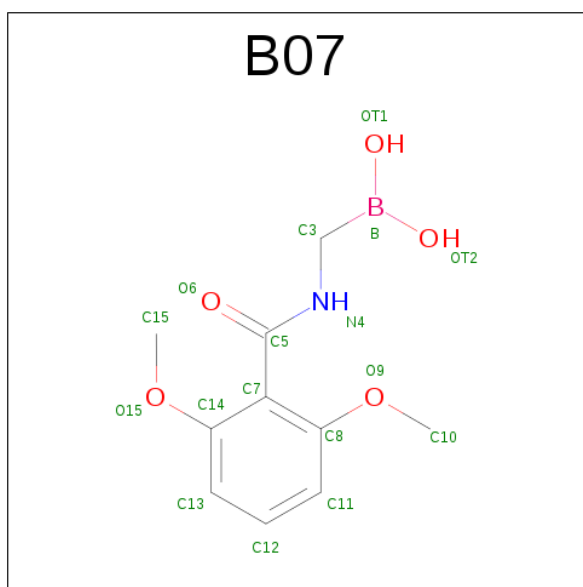
There are 4 unique types of molecules in this entry. The entry contains 13538 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 (three-letter code: B07) (formula: C₁₀H₁₄BNO₅).



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 (three-letter code: 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		

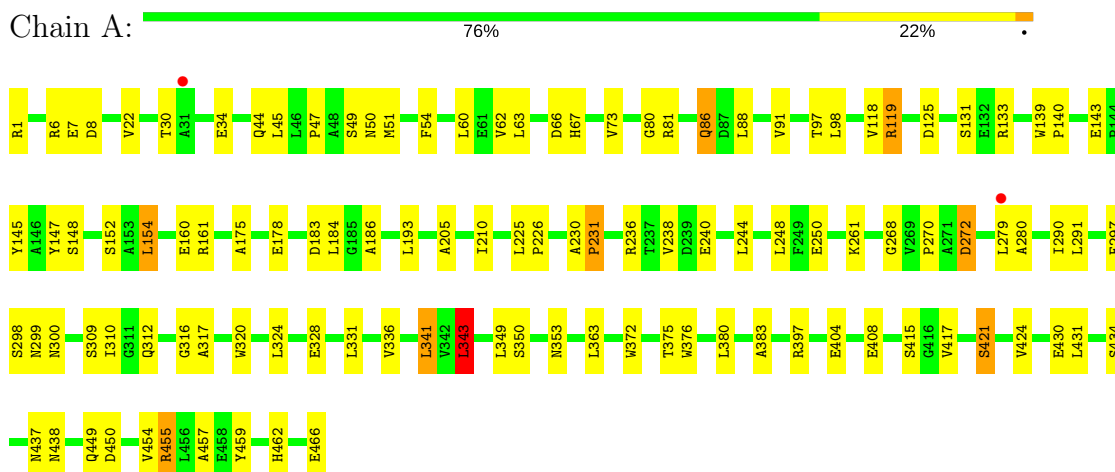
- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

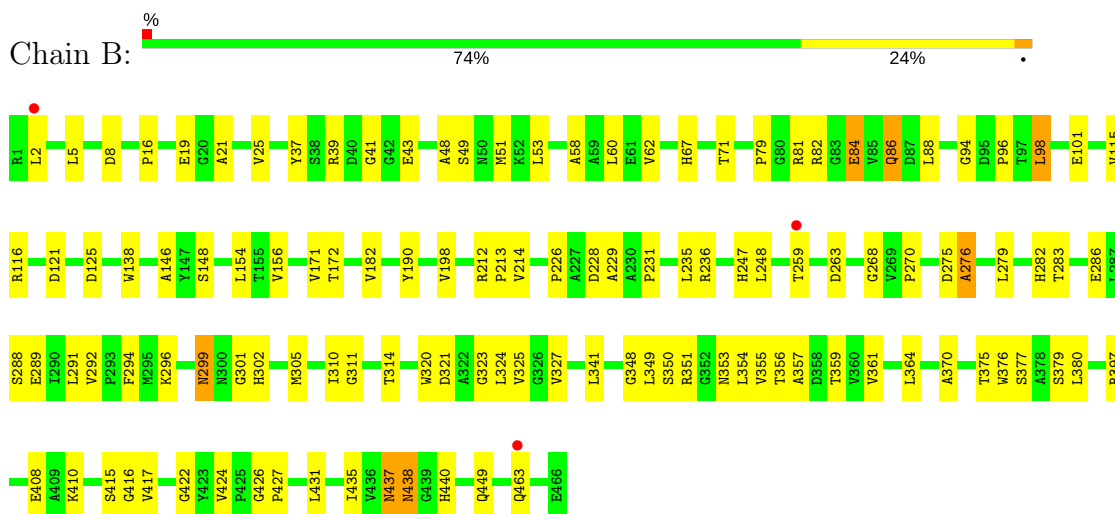
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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

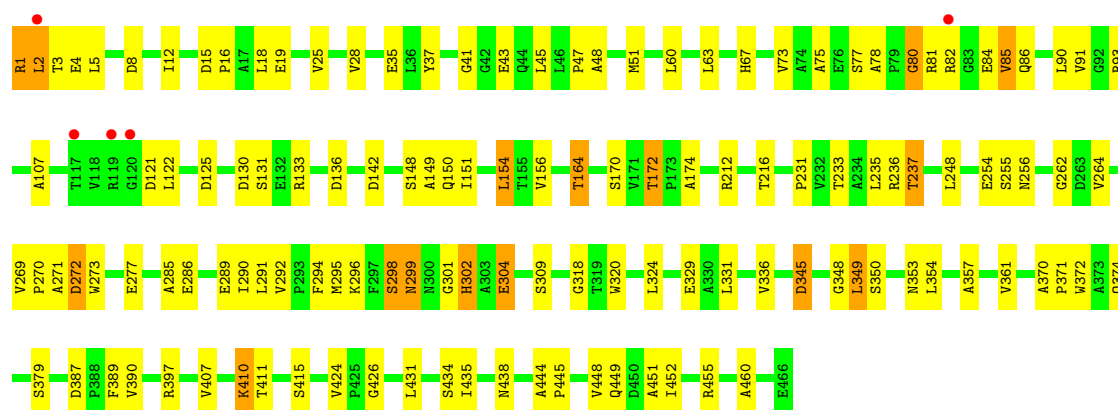


• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE

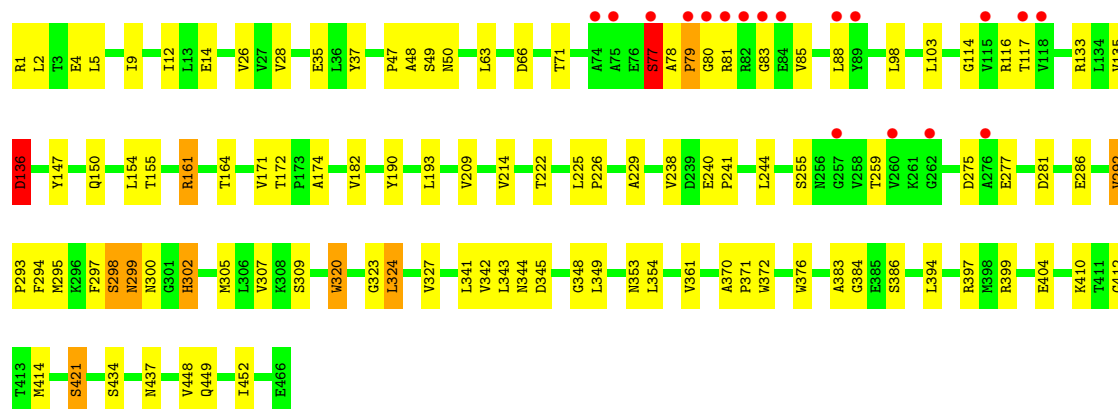
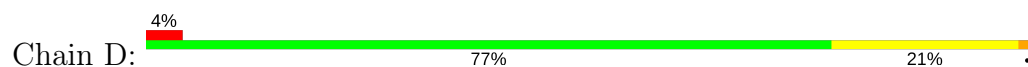


• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE





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



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) 95.5 (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	1770 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	59.9	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 26.5	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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, SO4, B07

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	0.46	0/3412	0.63	2/4666 (0.0%)
1	B	0.45	0/3403	0.62	0/4656
1	C	0.46	0/3403	0.62	0/4656
1	D	0.44	0/3412	0.60	1/4666 (0.0%)
All	All	0.45	0/13630	0.62	3/18644 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	154	LEU	CA-CB-CG	6.80	130.94	115.30
1	D	103	LEU	CA-CB-CG	6.37	129.95	115.30
1	A	343	LEU	CA-CB-CG	5.46	127.85	115.30

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	60	0
1	B	3344	0	3193	69	0
1	C	3344	0	3193	75	0
1	D	3353	0	3199	56	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	259	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 (259) 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: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: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:2:LEU:C	1:C:4:GLU:H	1.93	0.71
1:A:97:THR:HG21	1:A:290:ILE:HG12	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:C:270:PRO:C	1:C:272:ASP:H	1.97	0.67
1:D:421:SER:HB3	1:D:434:SER:HA	1.74	0.67
1:D:50:ASN:ND2	1:D:421:SER:OG	2.27	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
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: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:172:THR:HG23	1:C:231:PRO:HB3	1.84	0.59
1:C:25:VAL:HG22	1:C:435:ILE:HG23	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:77:SER:HA	1:D:277:GLU:HG3	1.87	0.56
1:B:51:MET:CE	1:B:353:ASN:HB3	2.36	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: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:D:150:GLN:HE22	1:D:240:GLU:H	1.54	0.55
1:D:49:SER:CB	1:D:412:GLY:HA2	2.32	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:2:LEU:C	1:C:4:GLU:N	2.63	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
1:B:437:ASN:C	1:B:438:ASN:HD22	2.14	0.51
1:C:142:ASP:O	1:C:148:SER:HB3	2.10	0.51
1:A:91:VAL:HG22	1:A:125:ASP:CB	2.38	0.51
1:B:299:ASN:HD22	1:B:299:ASN:C	2.14	0.50
1:B:41:GLY:HA2	1:B:357:ALA:HB3	1.94	0.50
1:B:212:ARG:O	1:B:213:PRO:C	2.50	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:270:PRO:C	1:C:272:ASP:N	2.64	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:A:51:MET:HE3	1:A:353:ASN:HB3	1.94	0.49
1:A:51:MET:HE2	1:A:353:ASN:HB3	1.95	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:135:VAL:O	1:D:136:ASP:C	2.51	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: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:96:PRO:HB2	1:B:302:HIS:CE1	2.49	0.48
1:B:94:GLY:O	1:B:96:PRO:HD3	2.13	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:448:VAL:O	1:C:452:ILE:HD12	2.13	0.47
1:C:90:LEU:HD23	1:C:154:LEU:HD11	1.95	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: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
1:A:383:ALA:O	1:A:404:GLU:HA	2.14	0.47
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: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:345:ASP:C	1:D:345:ASP:OD1	2.54	0.46
1:D:14:GLU:HA	1:D:14:GLU:OE1	2.16	0.46
1:D:384:GLY:HA3	1:D:404:GLU:HG3	1.97	0.46
1:A:270:PRO:C	1:A:272:ASP:H	2.19	0.46
1:B:275:ASP:O	1:B:276:ALA:C	2.54	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: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:161:ARG:HG2	1:D:297:PHE:CD2	2.52	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:63:LEU:O	1:A:67:HIS:HB2	2.17	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:PRO:HB2	1:B:302:HIS:ND1	2.33	0.44
1:A:186:ALA:HB1	1:A:248:LEU:HD11	1.99	0.44
1:A:45:LEU:HB3	1:A:438:ASN:OD1	2.18	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
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: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:370:ALA:C	1:D:372:TRP:H	2.20	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:133:ARG:HB3	1:C:150:GLN:HG2	2.01	0.43
1:C:73:VAL:HG22	1:C:90:LEU:HD12	2.00	0.43
1:C:170:SER:HB3	1:C:233:THR:HG23	2.00	0.43
1:C:78:ALA:N	1:C:277:GLU:HG2	2.33	0.43
1:D:292:VAL:HB	1:D:293:PRO:CD	2.49	0.43
1:B:323:GLY:O	1:B:327:VAL:HG23	2.19	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: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: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:372:TRP:O	1:C:372:TRP:HE3	2.02	0.42
1:C:45:LEU:HD22	1:C:438:ASN:HB2	1.99	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:A:341:LEU:HD22	1:A:343:LEU:HD22	2.00	0.42
1:A:73:VAL:HB	1:A:280:ALA:HB3	2.02	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:146:ALA:C	1:B:148:SER:H	2.23	0.42
1:B:341:LEU:CD2	1:B:355:VAL:HG12	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:341:LEU:HD13	1:D:343:LEU:HD11	2.01	0.42
1:D:12:ILE:HG21	1:D:452:ILE:HG12	2.02	0.42
1:B:296:LYS:NZ	1:B:380:LEU:O	2.51	0.41
1:B:311:GLY:CA	1:B:323:GLY:HA2	2.51	0.41
1:B:341:LEU:HG	1:B:359:THR:HG21	2.01	0.41
1:C:47:PRO:HG3	1:C:51:MET:HE3	2.01	0.41
1:A:47:PRO:HG3	1:A:51:MET:HE1	2.01	0.41
1:C:286:GLU:HB2	1:C:289:GLU:HG2	2.03	0.41
1:C:60:LEU:HD13	1:C:372:TRP:CE3	2.55	0.41
1:A:225:LEU:HA	1:A:226:PRO:HD3	1.91	0.41
1:B:51:MET:HE3	1:B:353:ASN:HB3	2.00	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:C:47:PRO:HG3	1:C:51:MET:HE2	2.03	0.40
1:A:240:GLU:HA	1:A:240:GLU:OE1	2.21	0.40
1:C:295:MET:HB2	1:C:379:SER:HB2	2.03	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:D:81:ARG:C	1:D:83:GLY:H	2.25	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: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
1:B:58:ALA:O	1:B:62:VAL:HG23	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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%)	28	67
1	B	464/466 (100%)	411 (89%)	48 (10%)	5 (1%)	17	54
1	C	464/466 (100%)	422 (91%)	34 (7%)	8 (2%)	11	42
1	D	464/466 (100%)	416 (90%)	42 (9%)	6 (1%)	14	48
All	All	1856/1864 (100%)	1667 (90%)	167 (9%)	22 (1%)	15	51

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%)	308 (91%)	31 (9%)	11	39
1	B	338/339 (100%)	316 (94%)	22 (6%)	20	55
1	C	338/339 (100%)	307 (91%)	31 (9%)	11	39
1	D	339/339 (100%)	304 (90%)	35 (10%)	8	32
All	All	1354/1356 (100%)	1235 (91%)	119 (9%)	12	42

All (119) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	7	GLU
1	A	22	VAL
1	A	34	GLU
1	A	44	GLN
1	A	66	ASP
1	A	81	ARG
1	A	98	LEU
1	A	119	ARG
1	A	154	LEU
1	A	160	GLU
1	A	183	ASP
1	A	250	GLU
1	A	272	ASP
1	A	279	LEU
1	A	291	LEU
1	A	299	ASN
1	A	309	SER
1	A	320	TRP
1	A	324	LEU
1	A	328	GLU
1	A	341	LEU
1	A	343	LEU
1	A	349	LEU
1	A	375	THR
1	A	376	TRP
1	A	415	SER
1	A	421	SER
1	A	437	ASN
1	A	455	ARG
1	A	466	GLU

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Mol	Chain	Res	Type
1	B	2	LEU
1	B	5	LEU
1	B	8	ASP
1	B	81	ARG
1	B	84	GLU
1	B	88	LEU
1	B	98	LEU
1	B	101	GLU
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	320	TRP
1	B	324	LEU
1	B	349	LEU
1	B	354	LEU
1	B	437	ASN
1	B	438	ASN
1	B	463	GLN
1	C	1	ARG
1	C	2	LEU
1	C	8	ASP
1	C	43	GLU
1	C	77	SER
1	C	84	GLU
1	C	85	VAL
1	C	93	ARG
1	C	136	ASP
1	C	154	LEU
1	C	164	THR
1	C	172	THR
1	C	237	THR
1	C	254	GLU
1	C	255	SER
1	C	269	VAL
1	C	272	ASP
1	C	291	LEU
1	C	298	SER
1	C	299	ASN

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Mol	Chain	Res	Type
1	C	302	HIS
1	C	304	GLU
1	C	309	SER
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	116	ARG
1	D	136	ASP
1	D	154	LEU
1	D	161	ARG
1	D	164	THR
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	255	SER
1	D	275	ASP
1	D	281	ASP
1	D	298	SER
1	D	299	ASN
1	D	302	HIS
1	D	305	MET
1	D	309	SER
1	D	320	TRP
1	D	324	LEU
1	D	344	ASN
1	D	349	LEU
1	D	376	TRP

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Mol	Chain	Res	Type
1	D	410	LYS
1	D	414	MET
1	D	421	SER
1	D	437	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	50	ASN
1	A	150	GLN
1	A	282	HIS
1	A	299	ASN
1	A	312	GLN
1	A	449	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	256	ASN
1	C	299	ASN
1	C	302	HIS
1	C	437	ASN
1	C	449	GLN
1	D	44	GLN
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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	$\# Z > 2$
2	B07	A	500	1	13,15,17	0.43	0	18,19,22	1.97	6 (33%)
3	SO4	A	601	-	4,4,4	0.16	0	6,6,6	0.21	0
3	SO4	A	602	-	4,4,4	0.15	0	6,6,6	0.12	0
3	SO4	A	603	-	4,4,4	0.14	0	6,6,6	0.17	0
3	SO4	A	604	-	4,4,4	0.18	0	6,6,6	0.26	0
2	B07	B	500	1	13,15,17	0.47	0	18,19,22	1.85	8 (44%)
3	SO4	B	601	-	4,4,4	0.18	0	6,6,6	0.15	0
3	SO4	B	602	-	4,4,4	0.17	0	6,6,6	0.07	0
3	SO4	B	603	-	4,4,4	0.17	0	6,6,6	0.19	0
3	SO4	B	604	-	4,4,4	0.13	0	6,6,6	0.15	0
2	B07	C	500	1	13,15,17	0.56	0	18,19,22	2.11	7 (38%)
3	SO4	C	601	-	4,4,4	0.21	0	6,6,6	0.13	0
3	SO4	C	602	-	4,4,4	0.20	0	6,6,6	0.09	0
3	SO4	C	603	-	4,4,4	0.21	0	6,6,6	0.21	0
3	SO4	C	604	-	4,4,4	0.15	0	6,6,6	0.14	0
2	B07	D	500	1	13,15,17	0.57	0	18,19,22	1.66	6 (33%)
3	SO4	D	601	-	4,4,4	0.23	0	6,6,6	0.16	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	D	602	-	4,4,4	0.15	0	6,6,6	0.15	0
3	SO4	D	603	-	4,4,4	0.18	0	6,6,6	0.21	0
3	SO4	D	604	-	4,4,4	0.15	0	6,6,6	0.11	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	-	0/10/11/13	0/1/1/1
3	SO4	A	601	-	-	0/0/0/0	0/0/0/0
3	SO4	A	602	-	-	0/0/0/0	0/0/0/0
3	SO4	A	603	-	-	0/0/0/0	0/0/0/0
3	SO4	A	604	-	-	0/0/0/0	0/0/0/0
2	B07	B	500	1	-	0/10/11/13	0/1/1/1
3	SO4	B	601	-	-	0/0/0/0	0/0/0/0
3	SO4	B	602	-	-	0/0/0/0	0/0/0/0
3	SO4	B	603	-	-	0/0/0/0	0/0/0/0
3	SO4	B	604	-	-	0/0/0/0	0/0/0/0
2	B07	C	500	1	-	0/10/11/13	0/1/1/1
3	SO4	C	601	-	-	0/0/0/0	0/0/0/0
3	SO4	C	602	-	-	0/0/0/0	0/0/0/0
3	SO4	C	603	-	-	0/0/0/0	0/0/0/0
3	SO4	C	604	-	-	0/0/0/0	0/0/0/0
2	B07	D	500	1	-	0/10/11/13	0/1/1/1
3	SO4	D	601	-	-	0/0/0/0	0/0/0/0
3	SO4	D	602	-	-	0/0/0/0	0/0/0/0
3	SO4	D	603	-	-	0/0/0/0	0/0/0/0
3	SO4	D	604	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	B07	O9-C8-C11	-3.41	118.64	124.37
2	D	500	B07	O15-C14-C13	-2.78	119.70	124.37
2	B	500	B07	O15-C14-C13	-2.73	119.79	124.37
2	D	500	B07	O9-C8-C11	-2.47	120.23	124.37
2	B	500	B07	O9-C8-C11	-2.41	120.32	124.37
2	C	500	B07	O15-C14-C13	-2.33	120.46	124.37
2	A	500	B07	C14-C7-C5	-2.23	116.79	121.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	B07	C3-N4-C5	2.09	125.69	121.89
2	C	500	B07	O15-C14-C7	2.13	118.98	115.84
2	D	500	B07	C10-O9-C8	2.16	120.64	117.54
2	A	500	B07	C15-O15-C14	2.22	120.73	117.54
2	C	500	B07	C8-C7-C14	2.27	121.36	117.87
2	D	500	B07	C7-C5-N4	2.29	119.79	115.51
2	C	500	B07	C7-C5-N4	2.43	120.05	115.51
2	B	500	B07	C7-C5-N4	2.48	120.16	115.51
2	B	500	B07	O9-C8-C7	2.49	119.51	115.84
2	B	500	B07	O15-C14-C7	2.70	119.81	115.84
2	B	500	B07	C10-O9-C8	2.73	121.47	117.54
2	D	500	B07	O15-C14-C7	2.77	119.92	115.84
2	A	500	B07	C10-O9-C8	2.93	121.76	117.54
2	B	500	B07	C15-O15-C14	2.96	121.79	117.54
2	D	500	B07	O9-C8-C7	3.02	120.28	115.84
2	C	500	B07	O9-C8-C7	3.09	120.39	115.84
2	A	500	B07	C7-C5-N4	3.14	121.40	115.51
2	C	500	B07	C15-O15-C14	3.85	123.07	117.54
2	A	500	B07	O9-C8-C7	3.88	121.55	115.84
2	C	500	B07	C10-O9-C8	4.89	124.58	117.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	B07	3	0
2	B	500	B07	1	0
2	C	500	B07	3	0
2	D	500	B07	4	0
3	D	603	SO4	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

6.1 Protein, DNA and RNA chains

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%)	-0.16	2 (0%) 92 84	35, 55, 84, 114	0
1	B	466/466 (100%)	-0.21	3 (0%) 89 77	33, 57, 90, 237	0
1	C	466/466 (100%)	-0.25	5 (1%) 80 65	37, 55, 83, 121	0
1	D	466/466 (100%)	-0.02	18 (3%) 40 19	44, 66, 107, 137	0
All	All	1864/1864 (100%)	-0.16	28 (1%) 74 54	33, 58, 96, 237	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	83	GLY	4.6
1	D	80	GLY	3.8
1	D	74	ALA	3.4
1	D	81	ARG	3.2
1	D	117	THR	3.2
1	B	2	LEU	3.1
1	C	117	THR	3.1
1	D	262	GLY	3.1
1	D	276	ALA	2.9
1	C	2	LEU	2.8
1	D	82	ARG	2.7
1	C	120	GLY	2.7
1	D	88	LEU	2.7
1	A	31	ALA	2.7
1	D	89	TYR	2.5
1	B	463	GLN	2.5
1	D	115	VAL	2.5
1	A	279	LEU	2.5
1	D	79	PRO	2.4
1	D	75	ALA	2.3
1	C	82	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	118	VAL	2.2
1	D	84	GLU	2.2
1	D	257	GLY	2.2
1	B	259	THR	2.1
1	D	260	VAL	2.1
1	D	77	SER	2.0
1	C	119	ARG	2.0

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	B	604	5/5	0.90	0.19	1.28	114,114,114,115	0
4	MG	A	611	1/1	0.95	0.19	0.54	64,64,64,64	0
2	B07	C	500	15/17	0.94	0.19	0.00	30,42,44,45	0
2	B07	D	500	15/17	0.96	0.19	-0.16	40,48,50,51	0
2	B07	A	500	15/17	0.96	0.20	-0.20	34,51,53,54	0
2	B07	B	500	15/17	0.96	0.21	-0.36	31,43,46,47	0
4	MG	A	610	1/1	0.87	0.08	-1.17	60,60,60,60	0
4	MG	D	611	1/1	0.98	0.10	-2.01	35,35,35,35	0
4	MG	D	610	1/1	0.95	0.06	-2.57	69,69,69,69	0
3	SO4	B	603	5/5	0.96	0.14	-	68,68,68,69	0
3	SO4	A	601	5/5	0.96	0.13	-	64,64,64,64	0
3	SO4	C	602	5/5	0.90	0.20	-	98,99,99,99	0
3	SO4	A	604	5/5	0.78	0.24	-	117,117,117,117	0
3	SO4	A	603	5/5	0.91	0.28	-	119,119,120,120	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	C	601	5/5	0.96	0.11	-	63,64,64,65	0
3	SO4	A	602	5/5	0.84	0.24	-	112,112,112,113	0
3	SO4	B	602	5/5	0.88	0.20	-	98,98,98,99	0
3	SO4	D	603	5/5	0.94	0.16	-	75,75,76,76	0
3	SO4	B	601	5/5	0.95	0.12	-	77,77,77,78	0
3	SO4	D	604	5/5	0.77	0.30	-	151,151,151,151	0
3	SO4	D	601	5/5	0.90	0.17	-	91,91,92,92	0
3	SO4	C	604	5/5	0.92	0.20	-	116,116,116,116	0
3	SO4	D	602	5/5	0.89	0.16	-	86,86,86,87	0
3	SO4	C	603	5/5	0.95	0.12	-	83,83,84,84	0

6.5 Other polymers [i](#)

There are no such residues in this entry.