



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 09:35 AM GMT

PDB ID : 2P83
Title : Potent and selective isophthalamide S2 hydroxyethylamine inhibitor of BACE1
Authors : Benson, T.E.; Prince, D.B.; Tomasselli, A.G.; Emmons, T.L.; Paddock, D.J.
Deposited on : 2007-03-21
Resolution : 2.50 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

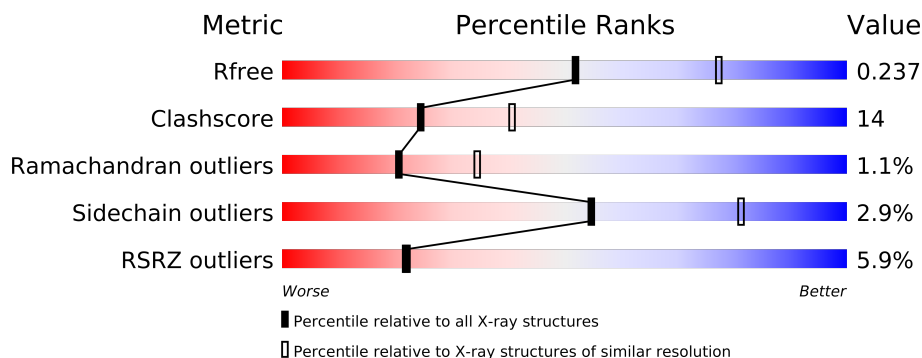
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.50 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}	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	455	
1	B	455	
1	C	455	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9104 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	371	Total	C	N	O	S	0	0	0
			2924	1874	487	549	14			
1	B	372	Total	C	N	O	S	0	0	0
			2929	1879	487	549	14			
1	C	373	Total	C	N	O	S	0	0	0
			2937	1883	488	552	14			

There are 45 discrepancies between the modelled and reference sequences:

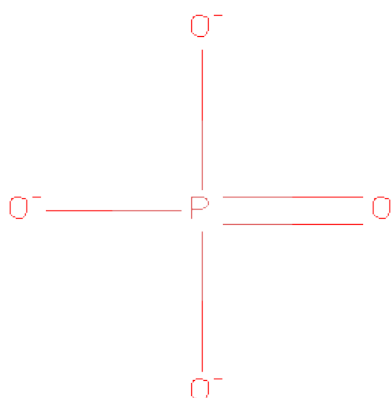
Chain	Residue	Modelled	Actual	Comment	Reference
A	1P	MET	-	EXPRESSION TAG	UNP P56817
A	2P	ALA	-	EXPRESSION TAG	UNP P56817
A	3P	SER	-	EXPRESSION TAG	UNP P56817
A	4P	MET	-	EXPRESSION TAG	UNP P56817
A	5P	THR	-	EXPRESSION TAG	UNP P56817
A	6P	GLY	-	EXPRESSION TAG	UNP P56817
A	7P	GLY	-	EXPRESSION TAG	UNP P56817
A	8P	GLN	-	EXPRESSION TAG	UNP P56817
A	9P	GLN	-	EXPRESSION TAG	UNP P56817
A	10P	MET	-	EXPRESSION TAG	UNP P56817
A	11P	GLY	-	EXPRESSION TAG	UNP P56817
A	12P	ARG	-	EXPRESSION TAG	UNP P56817
A	13P	GLY	-	EXPRESSION TAG	UNP P56817
A	14P	SER	-	EXPRESSION TAG	UNP P56817
A	15P	MET	-	EXPRESSION TAG	UNP P56817
B	1P	MET	-	EXPRESSION TAG	UNP P56817
B	2P	ALA	-	EXPRESSION TAG	UNP P56817
B	3P	SER	-	EXPRESSION TAG	UNP P56817
B	4P	MET	-	EXPRESSION TAG	UNP P56817
B	5P	THR	-	EXPRESSION TAG	UNP P56817
B	6P	GLY	-	EXPRESSION TAG	UNP P56817
B	7P	GLY	-	EXPRESSION TAG	UNP P56817
B	8P	GLN	-	EXPRESSION TAG	UNP P56817

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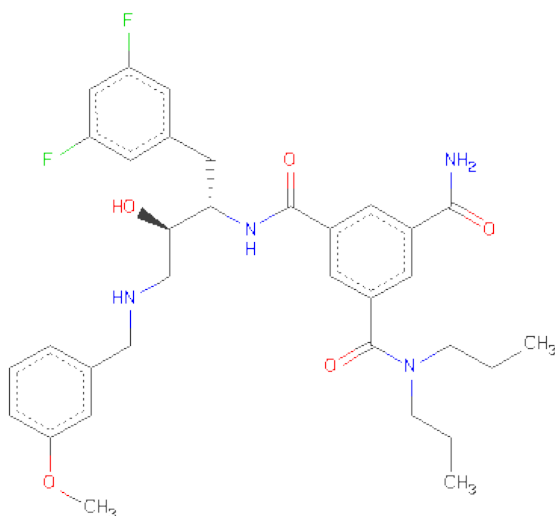
Chain	Residue	Modelled	Actual	Comment	Reference
B	9P	GLN	-	EXPRESSION TAG	UNP P56817
B	10P	MET	-	EXPRESSION TAG	UNP P56817
B	11P	GLY	-	EXPRESSION TAG	UNP P56817
B	12P	ARG	-	EXPRESSION TAG	UNP P56817
B	13P	GLY	-	EXPRESSION TAG	UNP P56817
B	14P	SER	-	EXPRESSION TAG	UNP P56817
B	15P	MET	-	EXPRESSION TAG	UNP P56817
C	1P	MET	-	EXPRESSION TAG	UNP P56817
C	2P	ALA	-	EXPRESSION TAG	UNP P56817
C	3P	SER	-	EXPRESSION TAG	UNP P56817
C	4P	MET	-	EXPRESSION TAG	UNP P56817
C	5P	THR	-	EXPRESSION TAG	UNP P56817
C	6P	GLY	-	EXPRESSION TAG	UNP P56817
C	7P	GLY	-	EXPRESSION TAG	UNP P56817
C	8P	GLN	-	EXPRESSION TAG	UNP P56817
C	9P	GLN	-	EXPRESSION TAG	UNP P56817
C	10P	MET	-	EXPRESSION TAG	UNP P56817
C	11P	GLY	-	EXPRESSION TAG	UNP P56817
C	12P	ARG	-	EXPRESSION TAG	UNP P56817
C	13P	GLY	-	EXPRESSION TAG	UNP P56817
C	14P	SER	-	EXPRESSION TAG	UNP P56817
C	15P	MET	-	EXPRESSION TAG	UNP P56817

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	C	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0

- Molecule 3 is N 3 -{(1S,2R)-1-(3,5-DIFLUOROBENZYL)-2-HYDROXY-3-[(3-METHOXY BENZYL)AMINO]PROPYL}-N 1 ,N 1 -DIPROPYLBENZENE-1,3,5-TRICARBOXAMID E (three-letter code: MR0) (formula: C₃₃H₄₀F₂N₄O₅).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C F N O 44 33 2 4 5	0	0
3	B	1	Total C F N O 44 33 2 4 5	0	0
3	C	1	Total C F N O 44 33 2 4 5	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	64	Total O 64 64	0	0
4	B	52	Total O 52 52	0	0

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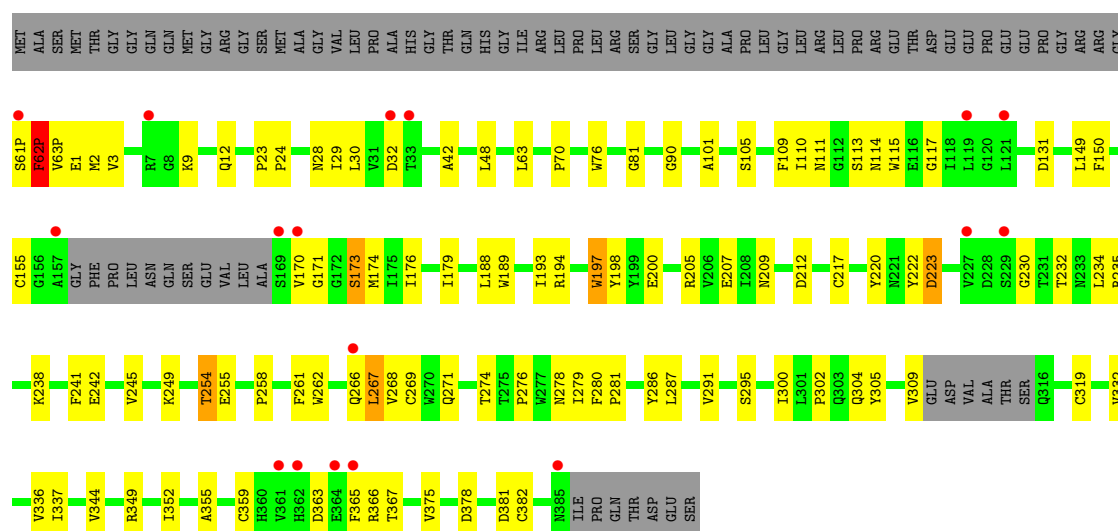
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	46	Total	O	0	0
			46	46		

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 errors 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.

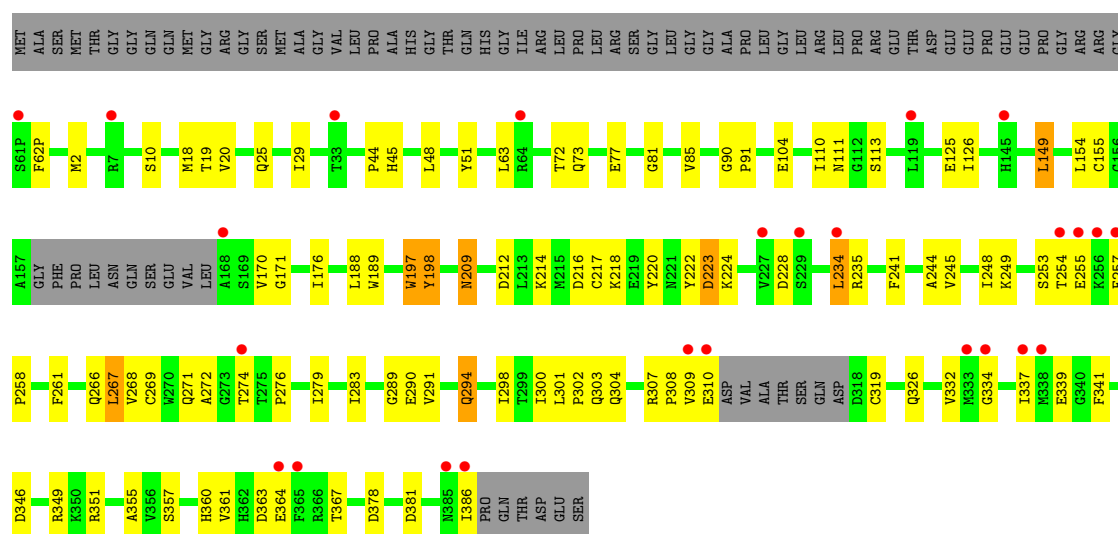
• Molecule 1: Beta-secretase 1

Chain A:



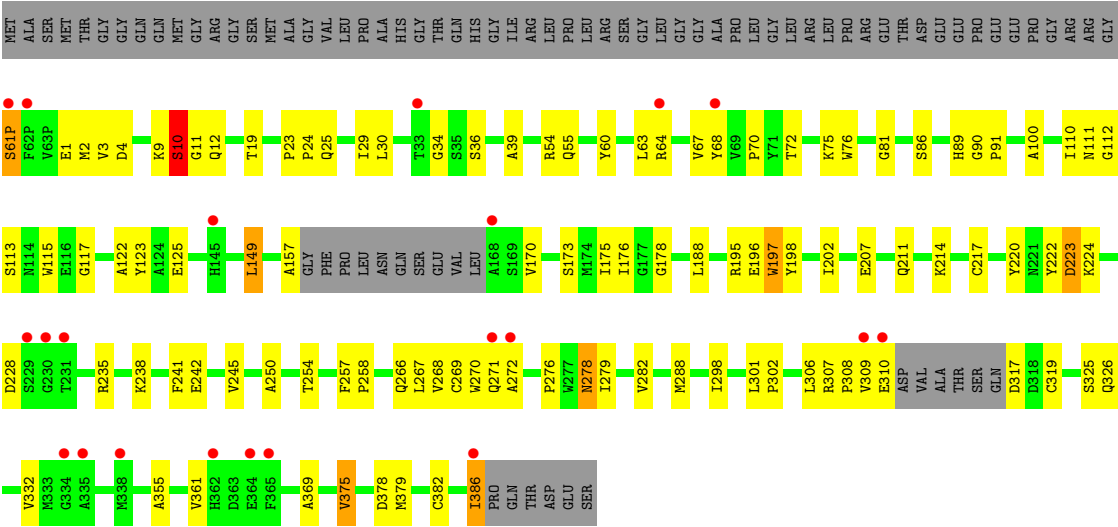
• Molecule 1: Beta-secretase 1

Chain B:



● Molecule 1: Beta-secretase 1

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.96Å 103.35Å 101.01Å 90.00° 104.12° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 31.61 – 2.47	Depositor EDS
% Data completeness (in resolution range)	83.7 (20.00-2.50) 82.4 (31.61-2.47)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.48Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.207 , 0.242 0.204 , 0.237	Depositor DCC
R_{free} test set	4859 reflections (11.21%)	DCC
Wilson B-factor (Å ²)	46.0	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 29.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 50612 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9104	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: PO4, MR0

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	0.39	0/2998	0.68	1/4072 (0.0%)
1	B	0.37	0/3003	0.68	1/4079 (0.0%)
1	C	0.37	0/3011	0.69	0/4090
All	All	0.38	0/9012	0.68	2/12241 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	193	ILE	N-CA-C	-5.12	97.18	111.00
1	B	198	TYR	N-CA-C	-5.02	97.43	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts i

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2924	0	2844	86	0
1	B	2929	0	2854	75	0
1	C	2937	0	2858	87	0
2	A	10	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	5	0	0	1	0
2	C	5	0	0	1	0
3	A	44	0	40	3	0
3	B	44	0	40	1	0
3	C	44	0	40	4	0
4	A	64	0	0	0	0
4	B	52	0	0	0	0
4	C	46	0	0	1	0
All	All	9104	0	8676	238	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (238) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:4:ASP:H	1:C:173:SER:HB3	1.27	0.96
1:C:269:CYS:HG	1:C:319:CYS:HG	0.88	0.87
1:C:4:ASP:N	1:C:173:SER:HB3	1.97	0.79
1:C:271:GLN:CD	1:C:271:GLN:H	1.91	0.74
1:A:9:LYS:HG3	1:A:12:GLN:HG3	1.69	0.74
1:A:375:VAL:CG2	1:C:375:VAL:HG22	2.18	0.73
1:A:149:LEU:HD23	1:A:150:PHE:N	2.02	0.73
1:B:276:PRO:O	1:B:279:ILE:HG12	1.89	0.73
1:A:375:VAL:HG22	1:C:375:VAL:HG22	1.73	0.71
1:C:188:LEU:HD23	1:C:355:ALA:HB2	1.73	0.70
1:C:202:ILE:CD1	1:C:379:MET:HG3	2.22	0.70
1:C:235:ARG:HG3	1:C:332:VAL:HB	1.72	0.69
1:B:291:VAL:HB	1:B:294:GLN:HG2	1.74	0.69
1:A:62(P):PHE:HB2	1:A:2:MET:CE	2.24	0.68
1:A:9:LYS:HD2	1:A:12:GLN:OE1	1.93	0.67
1:A:62(P):PHE:HB2	1:A:2:MET:HE2	1.77	0.67
1:C:271:GLN:CD	1:C:271:GLN:N	2.48	0.66
1:C:61(P):SER:N	1:C:175:ILE:CG2	2.60	0.65
1:B:155:CYS:O	1:B:170:VAL:HG13	1.97	0.65
1:B:267:LEU:H	1:B:267:LEU:HD12	1.62	0.64
1:B:241:PHE:O	1:B:245:VAL:HG23	1.98	0.64
1:A:261:PHE:CD1	1:A:268:VAL:HG23	2.33	0.63
1:B:291:VAL:O	1:B:294:GLN:HG3	1.98	0.63
1:A:300:ILE:HG12	1:A:337:ILE:CD1	2.28	0.63
1:C:61(P):SER:N	1:C:175:ILE:HG23	2.14	0.62
1:C:149:LEU:O	1:C:149:LEU:HD23	1.99	0.62
1:B:155:CYS:O	1:B:170:VAL:CG1	2.49	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:30:LEU:HD11	3:A:801:MR0:H121	1.82	0.60
1:C:68:TYR:OH	1:C:75:LYS:HD3	2.01	0.60
1:B:235:ARG:HG3	1:B:332:VAL:HB	1.84	0.60
1:B:244:ALA:O	1:B:248:ILE:HG13	2.01	0.60
1:A:188:LEU:HD23	1:A:355:ALA:HB2	1.84	0.59
1:A:235:ARG:HG3	1:A:332:VAL:HB	1.84	0.59
1:B:188:LEU:HD23	1:B:355:ALA:HB2	1.84	0.59
1:A:269:CYS:HG	1:A:319:CYS:CB	2.15	0.59
1:A:276:PRO:O	1:A:279:ILE:HG12	2.03	0.59
1:C:235:ARG:CG	1:C:332:VAL:HB	2.33	0.58
1:C:267:LEU:HD22	1:C:309:VAL:HG23	1.84	0.58
1:C:301:LEU:HB3	1:C:302:PRO:HD2	1.86	0.58
1:B:258:PRO:HG3	1:B:266:GLN:NE2	2.18	0.58
1:A:222:TYR:O	1:A:223:ASP:CB	2.51	0.57
1:C:288:MET:CE	1:C:378:ASP:HA	2.33	0.57
1:A:110:ILE:HB	1:A:113:SER:HB3	1.87	0.57
1:C:278:ASN:H	1:C:278:ASN:HD22	1.53	0.57
1:A:235:ARG:NH2	2:A:703:PO4:O2	2.37	0.57
1:B:245:VAL:HG12	1:B:249:LYS:HE3	1.87	0.57
1:B:216:ASP:OD1	1:B:218:LYS:HB2	2.03	0.57
1:A:62(P):PHE:CB	1:A:2:MET:HE2	2.35	0.56
1:C:157:ALA:HB2	1:C:170:VAL:HG12	1.87	0.56
1:A:375:VAL:HG23	1:C:375:VAL:HG22	1.87	0.56
1:B:301:LEU:HB3	1:B:302:PRO:HD2	1.87	0.56
1:A:155:CYS:O	1:A:170:VAL:CG1	2.53	0.56
1:C:19:THR:OG1	1:C:86:SER:HB2	2.05	0.56
1:A:375:VAL:HG22	1:C:375:VAL:HG13	1.88	0.55
1:C:238:LYS:HG3	1:C:326:GLN:OE1	2.05	0.55
1:A:378:ASP:HB3	1:A:381:ASP:OD2	2.06	0.55
1:A:3:VAL:O	1:A:173:SER:CB	2.54	0.55
1:A:249:LYS:HE2	1:A:262:TRP:CD1	2.42	0.55
1:A:61(P):SER:O	1:A:62(P):PHE:HB2	2.06	0.55
1:A:3:VAL:O	1:A:173:SER:HB3	2.07	0.55
1:C:19:THR:HA	1:C:25:GLN:O	2.07	0.54
1:B:212:ASP:O	1:B:214:LYS:HG3	2.07	0.54
1:B:170:VAL:HG12	1:B:171:GLY:N	2.22	0.54
1:A:62(P):PHE:CB	1:A:2:MET:CE	2.85	0.54
1:C:235:ARG:HB2	1:C:332:VAL:HB	1.90	0.54
1:B:253:SER:O	1:B:255:GLU:N	2.40	0.53
1:A:300:ILE:HG12	1:A:337:ILE:HD12	1.88	0.53
1:C:202:ILE:HD13	1:C:379:MET:HG3	1.91	0.53
1:B:2:MET:HG2	1:B:90:GLY:HA2	1.89	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:238:LYS:HE2	1:C:242:GLU:OE2	2.08	0.53
1:B:267:LEU:N	1:B:267:LEU:HD12	2.23	0.52
1:A:367:THR:H	1:C:211:GLN:HE22	1.58	0.52
1:B:269:CYS:HG	1:B:319:CYS:CB	2.20	0.52
1:C:241:PHE:CG	1:C:326:GLN:HB3	2.43	0.52
1:C:110:ILE:HB	1:C:113:SER:HB3	1.91	0.52
1:C:63:LEU:HG	1:C:81:GLY:HA2	1.92	0.52
1:C:125:GLU:CD	1:C:195:ARG:HH11	2.13	0.52
1:C:222:TYR:O	1:C:223:ASP:CB	2.57	0.52
1:A:287:LEU:O	1:A:295:SER:HB2	2.10	0.52
1:A:258:PRO:HG3	1:A:266:GLN:NE2	2.25	0.52
1:C:198:TYR:CE2	1:C:224:LYS:HE3	2.45	0.52
1:B:300:ILE:HA	1:B:304:GLN:OE1	2.10	0.52
1:B:303:GLN:HB2	1:B:361:VAL:CG1	2.39	0.51
1:B:125:GLU:HG2	1:B:197:TRP:HB3	1.92	0.51
1:A:3:VAL:O	1:A:173:SER:OG	2.28	0.51
1:C:270:TRP:O	1:C:317:ASP:HB3	2.11	0.51
1:B:378:ASP:HB3	1:B:381:ASP:OD2	2.11	0.51
1:B:228:ASP:O	1:B:334:GLY:HA2	2.11	0.50
1:A:42:ALA:CB	1:A:101:ALA:HB1	2.41	0.50
1:B:301:LEU:HD11	1:B:367:THR:HA	1.92	0.50
1:A:302:PRO:HA	1:A:305:TYR:CE2	2.47	0.50
1:C:34:GLY:O	3:C:803:MR0:H25	2.12	0.50
1:A:217:CYS:HG	1:A:382:CYS:HG	0.54	0.50
1:A:155:CYS:HG	1:A:359:CYS:HG	0.62	0.49
1:C:288:MET:HE3	1:C:378:ASP:HA	1.94	0.49
1:B:126:ILE:HG23	1:B:197:TRP:HB2	1.93	0.49
1:A:205:ARG:HB3	1:A:286:TYR:HB2	1.94	0.49
1:B:91:PRO:HD3	1:B:176:ILE:HB	1.95	0.49
1:C:29:ILE:HD12	1:C:117:GLY:HA3	1.94	0.49
1:B:72:THR:HB	3:B:802:MR0:H3	1.94	0.49
1:B:364:GLU:CD	1:B:364:GLU:N	2.66	0.49
1:B:290:GLU:N	1:B:294:GLN:OE1	2.37	0.49
1:A:291:VAL:CG1	1:B:189:TRP:CZ3	2.95	0.49
1:B:271:GLN:CD	1:B:271:GLN:H	2.16	0.49
1:B:289:GLY:HA3	1:B:294:GLN:O	2.12	0.49
1:A:194:ARG:HD3	1:A:200:GLU:OE2	2.13	0.49
1:A:278:ASN:OD1	1:C:254:THR:HG23	2.13	0.49
1:B:77:GLU:HG2	1:B:104:GLU:HB2	1.95	0.49
1:C:36:SER:OG	1:C:122:ALA:HB3	2.13	0.49
1:A:363:ASP:HB3	1:A:366:ARG:O	2.13	0.48
1:A:170:VAL:HG12	1:A:171:GLY:N	2.28	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:245:VAL:CG1	1:B:249:LYS:HE3	2.43	0.48
1:A:235:ARG:HB2	1:A:332:VAL:HB	1.95	0.48
1:A:63:LEU:HG	1:A:81:GLY:HA2	1.95	0.48
1:B:198:TYR:CE2	1:B:224:LYS:HE3	2.49	0.48
1:C:2:MET:HG2	1:C:90:GLY:HA2	1.95	0.48
1:B:267:LEU:HD23	1:B:309:VAL:HG21	1.95	0.48
1:B:217:CYS:HA	1:B:220:TYR:CD1	2.48	0.48
1:B:302:PRO:HG2	1:B:303:GLN:OE1	2.13	0.48
1:C:72:THR:HB	3:C:803:MR0:H3	1.95	0.48
1:C:302:PRO:O	1:C:306:LEU:HB2	2.14	0.48
1:A:359:CYS:O	1:A:359:CYS:SG	2.72	0.47
1:A:365:PHE:CG	1:C:250:ALA:HB1	2.49	0.47
1:C:307:ARG:HA	1:C:308:PRO:HD3	1.72	0.47
1:A:63(P):VAL:HG13	1:A:1:GLU:N	2.28	0.47
1:B:222:TYR:O	1:B:223:ASP:HB3	2.14	0.47
1:C:235:ARG:NH2	2:C:701:PO4:O2	2.30	0.47
1:B:20:VAL:HG12	1:B:85:VAL:HG22	1.96	0.47
1:A:300:ILE:HG23	1:A:304:GLN:HB2	1.97	0.47
1:A:110:ILE:HD11	3:A:801:MR0:H101	1.97	0.47
1:C:125:GLU:O	1:C:125:GLU:HG3	2.13	0.47
1:C:298:ILE:HG13	1:C:298:ILE:O	2.15	0.47
1:B:149:LEU:HD23	1:B:149:LEU:C	2.34	0.47
1:A:9:LYS:HG3	1:A:12:GLN:CG	2.42	0.47
1:B:44:PRO:HD3	1:B:51:TYR:CZ	2.50	0.47
1:B:298:ILE:HG13	1:B:298:ILE:O	2.15	0.47
1:C:267:LEU:O	1:C:267:LEU:HD12	2.14	0.47
1:C:10:SER:HB3	1:C:11:GLY:H	1.49	0.47
1:B:19:THR:HA	1:B:25:GLN:O	2.15	0.47
1:C:197:TRP:CG	1:C:198:TYR:N	2.84	0.46
1:A:149:LEU:C	1:A:149:LEU:HD23	2.35	0.46
1:A:170:VAL:HG12	1:A:171:GLY:H	1.79	0.46
1:C:12:GLN:OE1	1:C:113:SER:HA	2.15	0.46
1:A:179:ILE:HD11	1:A:344:VAL:HG11	1.97	0.46
1:B:357:SER:O	1:B:360:HIS:HB3	2.16	0.46
1:B:222:TYR:O	1:B:223:ASP:CB	2.63	0.46
1:B:257:PHE:HD2	1:B:268:VAL:HG21	1.80	0.46
1:B:267:LEU:HD23	1:B:309:VAL:CG2	2.46	0.46
1:A:207:GLU:HG2	1:A:212:ASP:HA	1.97	0.46
1:C:238:LYS:O	1:C:242:GLU:HG3	2.15	0.46
1:B:283:ILE:HB	1:B:300:ILE:HG12	1.98	0.46
1:B:110:ILE:HB	1:B:113:SER:HB3	1.97	0.46
1:B:234:LEU:HG	1:B:337:ILE:HD11	1.98	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:282:VAL:HG12	1:C:301:LEU:HD23	1.98	0.45
1:B:386:ILE:H	1:B:386:ILE:HG13	1.44	0.45
1:C:217:CYS:HA	1:C:220:TYR:CD1	2.51	0.45
1:C:9:LYS:HE2	1:C:12:GLN:OE1	2.16	0.45
1:A:217:CYS:HA	1:A:220:TYR:CD1	2.51	0.45
1:C:149:LEU:HD21	1:C:178:GLY:HA2	1.97	0.45
1:A:344:VAL:O	1:A:352:ILE:HA	2.17	0.45
1:C:241:PHE:O	1:C:245:VAL:HG23	2.17	0.45
1:B:209:ASN:HD22	1:B:209:ASN:HA	1.62	0.45
1:B:303:GLN:HB2	1:B:361:VAL:HG11	1.98	0.45
1:A:29:ILE:HD12	1:A:117:GLY:HA3	1.98	0.45
1:C:110:ILE:HD11	3:C:803:MR0:H101	1.98	0.45
1:B:271:GLN:O	1:B:272:ALA:C	2.54	0.45
1:B:346:ASP:HB3	1:B:351:ARG:HG3	1.98	0.45
1:A:189:TRP:CZ3	1:B:291:VAL:CG1	3.00	0.44
1:B:261:PHE:CD1	1:B:268:VAL:HG23	2.52	0.44
1:C:235:ARG:CB	1:C:332:VAL:HB	2.48	0.44
1:A:267:LEU:HD23	1:A:309:VAL:HG21	2.00	0.44
1:B:294:GLN:HE21	1:B:294:GLN:HB2	1.63	0.44
1:C:369:ALA:HB1	4:C:952:HOH:O	2.18	0.44
1:A:28:ASN:HB2	1:A:115:TRP:HA	1.98	0.44
1:C:214:LYS:HB3	1:C:214:LYS:HE2	1.81	0.44
1:A:155:CYS:CB	1:A:359:CYS:SG	3.05	0.44
1:C:30:LEU:HD23	1:C:30:LEU:C	2.38	0.44
1:A:174:MET:CE	1:A:176:ILE:HD11	2.48	0.44
1:C:207:GLU:HA	1:C:211:GLN:O	2.18	0.44
1:C:34:GLY:HA3	1:C:228:ASP:OD1	2.17	0.44
1:A:113:SER:O	1:A:114:ASN:HB3	2.17	0.44
1:A:30:LEU:CD2	1:A:32:ASP:HB2	2.48	0.43
1:A:197:TRP:CG	1:A:198:TYR:N	2.86	0.43
1:B:271:GLN:HB2	1:B:274:THR:HG21	2.01	0.43
1:A:238:LYS:O	1:A:242:GLU:HG3	2.18	0.43
1:B:170:VAL:CG1	1:B:171:GLY:N	2.81	0.43
1:A:76:TRP:HA	1:A:105:SER:HA	1.99	0.43
1:C:54:ARG:HD2	1:C:60:TYR:CZ	2.53	0.43
1:C:257:PHE:HD2	1:C:268:VAL:HG21	1.82	0.43
1:A:268:VAL:HG12	1:A:269:CYS:N	2.34	0.43
1:C:91:PRO:HD3	1:C:176:ILE:HB	2.00	0.43
1:A:230:GLY:O	3:A:801:MR0:H122	2.19	0.43
1:A:280:PHE:HB3	1:A:302:PRO:HB3	2.00	0.43
1:C:39:ALA:HB2	1:C:100:ALA:HB3	2.01	0.43
1:B:63:LEU:HG	1:B:81:GLY:HA2	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:45:HIS:HB3	1:B:48:LEU:HG	2.01	0.43
1:C:67:VAL:HG13	1:C:76:TRP:HZ3	1.84	0.43
1:B:268:VAL:HG12	1:B:269:CYS:N	2.33	0.42
1:A:48:LEU:HD21	1:A:109:PHE:CE2	2.53	0.42
1:C:217:CYS:HG	1:C:382:CYS:CB	2.26	0.42
1:C:9:LYS:O	1:C:10:SER:C	2.57	0.42
1:A:254:THR:HG22	1:A:255:GLU:HG3	2.01	0.42
1:A:23:PRO:HA	1:A:24:PRO:HD3	1.82	0.42
1:C:12:GLN:C	3:C:803:MR0:H111	2.40	0.42
1:B:154:LEU:O	1:B:339:GLU:HA	2.19	0.42
1:C:267:LEU:HD12	1:C:267:LEU:C	2.40	0.42
1:C:23:PRO:HA	1:C:24:PRO:HD3	1.88	0.42
1:C:30:LEU:HD12	1:C:115:TRP:CE2	2.54	0.42
1:B:18:MET:SD	1:B:29:ILE:HG13	2.60	0.42
1:C:3:VAL:O	1:C:4:ASP:HB3	2.20	0.42
1:A:271:GLN:O	1:A:274:THR:HG23	2.20	0.42
1:A:2:MET:HG2	1:A:90:GLY:HA2	2.02	0.41
1:B:241:PHE:CG	1:B:326:GLN:HB3	2.55	0.41
1:C:149:LEU:C	1:C:149:LEU:HD23	2.39	0.41
1:C:386:ILE:O	1:C:386:ILE:HG22	2.20	0.41
1:A:365:PHE:CD1	1:A:366:ARG:HG3	2.54	0.41
1:C:276:PRO:O	1:C:279:ILE:HG12	2.20	0.41
1:C:2:MET:HG2	1:C:89:HIS:O	2.19	0.41
1:A:174:MET:HE3	1:A:176:ILE:HD11	2.01	0.41
1:A:62(P):PHE:CB	1:A:2:MET:HE3	2.49	0.41
1:B:235:ARG:NH2	2:B:704:PO4:O1	2.53	0.41
1:A:48:LEU:HD21	1:A:109:PHE:CD2	2.55	0.41
1:B:307:ARG:HA	1:B:308:PRO:HD3	1.74	0.41
1:A:349:ARG:HH21	1:B:349:ARG:HH21	1.67	0.41
1:B:309:VAL:O	1:B:310:GLU:C	2.59	0.41
1:B:341:PHE:HB3	1:B:355:ALA:O	2.21	0.41
1:B:303:GLN:NE2	1:B:363:ASP:HB3	2.35	0.41
1:C:123:TYR:CZ	1:C:196:GLU:HG2	2.54	0.41
1:A:209:ASN:ND2	1:A:281:PRO:HB3	2.36	0.41
1:A:235:ARG:CG	1:A:332:VAL:HB	2.50	0.41
1:A:241:PHE:O	1:A:245:VAL:HG23	2.21	0.41
1:C:258:PRO:HB2	1:C:266:GLN:OE1	2.21	0.41
1:A:304:GLN:O	1:A:336:VAL:HB	2.21	0.40
1:A:232:THR:O	1:A:336:VAL:HG13	2.20	0.40
1:A:268:VAL:O	1:A:319:CYS:HA	2.22	0.40
1:C:241:PHE:CD2	1:C:326:GLN:HB3	2.56	0.40
1:C:9:LYS:HE2	1:C:112:GLY:O	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:310:GLU:HG3	1:C:310:GLU:O	2.21	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	365/455 (80%)	340 (93%)	21 (6%)	4 (1%)	21	34
1	B	366/455 (80%)	336 (92%)	26 (7%)	4 (1%)	21	34
1	C	367/455 (81%)	345 (94%)	18 (5%)	4 (1%)	21	34
All	All	1098/1365 (80%)	1021 (93%)	65 (6%)	12 (1%)	21	34

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62(P)	PHE
1	A	223	ASP
1	B	10	SER
1	B	223	ASP
1	C	10	SER
1	C	223	ASP
1	B	254	THR
1	C	272	ALA
1	B	73	GLN
1	A	131	ASP
1	A	70	PRO
1	C	70	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	317/381 (83%)	310 (98%)	7 (2%)	64	88
1	B	317/381 (83%)	309 (98%)	8 (2%)	60	85
1	C	318/381 (84%)	305 (96%)	13 (4%)	41	67
All	All	952/1143 (83%)	924 (97%)	28 (3%)	55	81

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62(P)	PHE
1	A	111	ASN
1	A	173	SER
1	A	197	TRP
1	A	234	LEU
1	A	254	THR
1	A	267	LEU
1	B	62(P)	PHE
1	B	111	ASN
1	B	149	LEU
1	B	197	TRP
1	B	209	ASN
1	B	234	LEU
1	B	267	LEU
1	B	294	GLN
1	C	61(P)	SER
1	C	1	GLU
1	C	10	SER
1	C	55	GLN
1	C	64	ARG
1	C	111	ASN
1	C	149	LEU
1	C	197	TRP
1	C	278	ASN
1	C	325	SER
1	C	361	VAL
1	C	375	VAL
1	C	386	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	98	ASN
1	A	111	ASN
1	A	266	GLN
1	A	271	GLN
1	B	73	GLN
1	B	92	ASN
1	B	98	ASN
1	B	111	ASN
1	B	209	ASN
1	B	266	GLN
1	C	98	ASN
1	C	111	ASN
1	C	211	GLN
1	C	278	ASN
1	C	294	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

7 ligands are modelled in this entry.

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	PO4	A	702	-	4,4,4	1.11	0	6,6,6	0.31	0
2	PO4	A	703	-	4,4,4	0.64	0	6,6,6	0.31	0
3	MR0	A	801	-	46,46,46	2.55	25 (54%)	62,62,62	1.69	13 (20%)
2	PO4	B	704	-	4,4,4	0.67	0	6,6,6	0.31	0
3	MR0	B	802	-	46,46,46	2.66	24 (52%)	62,62,62	1.79	15 (24%)
2	PO4	C	701	-	4,4,4	1.02	0	6,6,6	0.31	0
3	MR0	C	803	-	46,46,46	2.49	25 (54%)	62,62,62	1.70	14 (22%)

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	PO4	A	702	-	-	0/0/0/0	0/0/0/0
2	PO4	A	703	-	-	0/0/0/0	0/0/0/0
3	MR0	A	801	-	-	0/42/42/42	0/3/3/3
2	PO4	B	704	-	-	0/0/0/0	0/0/0/0
3	MR0	B	802	-	-	0/42/42/42	0/3/3/3
2	PO4	C	701	-	-	0/0/0/0	0/0/0/0
3	MR0	C	803	-	-	0/42/42/42	0/3/3/3

All (74) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	802	MR0	C25-C20	5.55	1.49	1.39
3	B	802	MR0	C29-C30	4.97	1.46	1.37
3	B	802	MR0	C29-C28	4.92	1.48	1.39
3	A	801	MR0	C5-C	4.80	1.46	1.39
3	C	803	MR0	C5-C	4.61	1.46	1.39
3	A	801	MR0	C25-C20	4.57	1.47	1.39
3	A	801	MR0	C29-C30	4.50	1.45	1.37
3	A	801	MR0	C31-C30	4.50	1.45	1.37
3	B	802	MR0	C5-C	4.49	1.46	1.39
3	B	802	MR0	C31-C30	4.47	1.45	1.37
3	C	803	MR0	C31-C30	4.42	1.45	1.37
3	C	803	MR0	C25-C20	4.34	1.47	1.39
3	C	803	MR0	C29-C30	4.29	1.45	1.37
3	B	802	MR0	C1-C2	4.29	1.45	1.39
3	B	802	MR0	C6-N	4.28	1.45	1.34
3	C	803	MR0	C6-N	4.19	1.45	1.34
3	A	801	MR0	C1-C2	4.18	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	MR0	C6-N	4.12	1.45	1.34
3	A	801	MR0	C29-C28	4.01	1.46	1.39
3	C	803	MR0	C29-C28	3.97	1.46	1.39
3	B	802	MR0	C8-N1	3.76	1.42	1.34
3	B	802	MR0	C3-C2	3.76	1.45	1.39
3	C	803	MR0	C1-C2	3.70	1.45	1.39
3	C	803	MR0	C5-C4	3.62	1.44	1.39
3	A	801	MR0	C27-C26	3.41	1.43	1.37
3	A	801	MR0	C5-C4	3.32	1.44	1.39
3	C	803	MR0	C31-C26	3.30	1.43	1.37
3	A	801	MR0	C31-C26	3.27	1.43	1.37
3	B	802	MR0	C31-C26	3.23	1.43	1.37
3	A	801	MR0	C7-N7A	3.21	1.40	1.33
3	B	802	MR0	C5-C4	3.18	1.44	1.39
3	A	801	MR0	C23-C24	3.12	1.45	1.38
3	A	801	MR0	C3-C2	3.08	1.44	1.39
3	C	803	MR0	C27-C26	3.05	1.43	1.37
3	C	803	MR0	C22-C23	3.05	1.46	1.39
3	C	803	MR0	C3-C2	3.02	1.44	1.39
3	C	803	MR0	C23-C24	2.99	1.44	1.38
3	A	801	MR0	C15-N1	2.93	1.51	1.46
3	C	803	MR0	C8-N1	2.91	1.40	1.34
3	B	802	MR0	C27-C26	2.87	1.42	1.37
3	B	802	MR0	C25-C24	2.78	1.44	1.38
3	B	802	MR0	C21-C20	2.77	1.45	1.38
3	A	801	MR0	C22-C21	2.76	1.45	1.39
3	B	802	MR0	C17-C28	2.75	1.58	1.51
3	C	803	MR0	C7-N7A	2.74	1.39	1.33
3	A	801	MR0	C22-C23	2.74	1.45	1.39
3	A	801	MR0	C8-N1	2.73	1.40	1.34
3	A	801	MR0	C1-C	2.66	1.43	1.39
3	C	803	MR0	C27-C28	2.63	1.44	1.39
3	B	802	MR0	C1-C	2.61	1.43	1.39
3	B	802	MR0	C22-C23	2.61	1.45	1.39
3	B	802	MR0	C15-N1	2.58	1.51	1.46
3	C	803	MR0	C15-N1	2.55	1.51	1.46
3	A	801	MR0	C27-C28	2.54	1.44	1.39
3	B	802	MR0	F26-C26	-2.53	1.30	1.36
3	A	801	MR0	C10-N	2.53	1.54	1.47
3	B	802	MR0	C7-N7A	2.51	1.38	1.33
3	A	801	MR0	C21-C20	2.49	1.44	1.38
3	C	803	MR0	C22-C21	2.47	1.44	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	803	MR0	C25-C24	2.47	1.43	1.38
3	B	802	MR0	C10-N	2.39	1.53	1.47
3	C	803	MR0	C2-C7	-2.38	1.46	1.50
3	B	802	MR0	C22-C21	2.33	1.44	1.39
3	A	801	MR0	C25-C24	2.28	1.43	1.38
3	B	802	MR0	C23-C24	2.27	1.43	1.38
3	C	803	MR0	F30-C30	-2.21	1.31	1.36
3	C	803	MR0	C10-N	2.18	1.53	1.47
3	A	801	MR0	C17-C28	2.17	1.56	1.51
3	C	803	MR0	C17-C28	2.16	1.56	1.51
3	A	801	MR0	F30-C30	-2.16	1.31	1.36
3	B	802	MR0	C19-C20	2.03	1.56	1.51
3	C	803	MR0	C21-C20	2.03	1.43	1.38
3	A	801	MR0	C18-N2	2.03	1.50	1.47
3	C	803	MR0	O-C6	2.01	1.26	1.22

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	803	MR0	C17-C15-N1	-5.07	104.28	110.20
3	A	801	MR0	C-C6-N	4.81	126.09	118.88
3	B	802	MR0	C17-C15-N1	-4.67	104.76	110.20
3	A	801	MR0	C17-C15-N1	-4.66	104.76	110.20
3	B	802	MR0	C-C6-N	4.63	125.83	118.88
3	C	803	MR0	C16-C15-N1	4.51	117.64	109.61
3	A	801	MR0	C16-C15-N1	4.37	117.39	109.61
3	C	803	MR0	C-C6-N	3.92	124.77	118.88
3	C	803	MR0	C28-C27-C26	3.85	122.14	118.96
3	B	802	MR0	C28-C27-C26	3.65	121.98	118.96
3	B	802	MR0	CO-O3-C24	3.53	126.14	117.54
3	A	801	MR0	C28-C27-C26	3.53	121.88	118.96
3	B	802	MR0	C31-C30-C29	-3.49	118.55	123.48
3	B	802	MR0	C20-C19-N2	-3.47	103.53	113.02
3	B	802	MR0	C16-C15-N1	3.34	115.56	109.61
3	C	803	MR0	C24-C25-C20	3.27	123.11	119.73
3	A	801	MR0	C20-C19-N2	-2.99	104.83	113.02
3	C	803	MR0	C31-C26-C27	-2.95	119.30	123.48
3	C	803	MR0	C30-C31-C26	2.88	120.66	116.24
3	A	801	MR0	C24-C25-C20	2.81	122.64	119.73
3	C	803	MR0	C20-C19-N2	-2.81	105.34	113.02
3	C	803	MR0	C31-C30-C29	-2.78	119.55	123.48
3	B	802	MR0	C30-C31-C26	2.77	120.49	116.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	MR0	C31-C26-C27	-2.76	119.57	123.48
3	A	801	MR0	C30-C31-C26	2.75	120.45	116.24
3	B	802	MR0	O-C6-C	-2.72	115.04	120.17
3	A	801	MR0	C31-C30-C29	-2.72	119.64	123.48
3	B	802	MR0	C23-C24-C25	-2.62	116.80	120.56
3	B	802	MR0	C16-C18-N2	-2.59	107.79	113.03
3	C	803	MR0	C23-C24-C25	-2.58	116.85	120.56
3	B	802	MR0	C24-C25-C20	2.50	122.31	119.73
3	A	801	MR0	O-C6-N	-2.41	117.85	122.33
3	B	802	MR0	O3-C24-C25	2.36	131.30	119.90
3	A	801	MR0	CO-O3-C24	2.27	123.08	117.54
3	A	801	MR0	C23-C24-C25	-2.21	117.38	120.56
3	A	801	MR0	O-C6-C	-2.18	116.06	120.17
3	B	802	MR0	C31-C26-C27	-2.11	120.49	123.48
3	B	802	MR0	C22-C23-C24	2.07	122.59	118.91
3	C	803	MR0	O-C6-C	-2.03	116.34	120.17
3	C	803	MR0	C3-C2-C1	2.03	122.16	119.64
3	C	803	MR0	CO-O3-C24	2.01	122.44	117.54
3	C	803	MR0	C16-C18-N2	-2.01	108.96	113.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	371/455 (81%)	0.12	17 (4%) 31 32	26, 41, 65, 81	0
1	B	372/455 (81%)	0.34	25 (6%) 17 17	23, 46, 74, 89	0
1	C	373/455 (81%)	0.23	21 (5%) 24 24	26, 44, 70, 86	0
All	All	1116/1365 (81%)	0.23	63 (5%) 22 24	23, 44, 71, 89	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	168	ALA	5.4
1	C	386	ILE	5.4
1	A	157	ALA	4.8
1	C	272	ALA	4.7
1	B	256	LYS	4.4
1	B	61(P)	SER	4.4
1	C	168	ALA	4.4
1	B	310	GLU	4.4
1	C	364	GLU	3.9
1	B	386	ILE	3.9
1	A	61(P)	SER	3.8
1	C	365	PHE	3.6
1	C	62(P)	PHE	3.3
1	B	254	THR	3.3
1	B	337	ILE	3.3
1	A	362	HIS	3.2
1	B	64	ARG	3.1
1	B	365	PHE	3.1
1	A	227	VAL	3.1
1	C	145	HIS	3.1
1	A	169	SER	3.0
1	C	61(P)	SER	3.0
1	B	309	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	7	ARG	2.9
1	A	365	PHE	2.8
1	C	33	THR	2.8
1	B	227	VAL	2.8
1	B	274	THR	2.7
1	C	310	GLU	2.7
1	C	229	SER	2.7
1	B	385	ASN	2.7
1	A	364	GLU	2.6
1	B	33	THR	2.6
1	B	364	GLU	2.5
1	A	119	LEU	2.5
1	B	234	LEU	2.5
1	B	145	HIS	2.4
1	B	229	SER	2.4
1	C	335	ALA	2.4
1	B	255	GLU	2.4
1	C	362	HIS	2.4
1	A	229	SER	2.4
1	A	170	VAL	2.3
1	B	334	GLY	2.3
1	C	334	GLY	2.3
1	B	7	ARG	2.3
1	B	257	PHE	2.3
1	A	266	GLN	2.2
1	C	309	VAL	2.2
1	A	121	LEU	2.2
1	C	64	ARG	2.2
1	B	119	LEU	2.2
1	B	333	MET	2.2
1	C	338	MET	2.2
1	C	231	THR	2.1
1	A	361	VAL	2.1
1	A	32	ASP	2.1
1	A	33	THR	2.1
1	C	68	TYR	2.1
1	C	230	GLY	2.0
1	C	271	GLN	2.0
1	A	385	ASN	2.0
1	B	338	MET	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

There are no carbohydrates in this entry.

6.4 Ligands

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MR0	C	803	44/44	0.21	-	36,40,45,49	0
2	PO4	B	704	5/5	0.41	-	99,99,99,99	0
2	PO4	A	703	5/5	0.34	-	96,96,97,98	0
2	PO4	A	702	5/5	0.41	-	99,99,99,99	0
3	MR0	A	801	44/44	0.18	-	29,35,39,43	0
3	MR0	B	802	44/44	0.21	-	38,43,45,49	0
2	PO4	C	701	5/5	0.43	-	99,99,99,99	0

6.5 Other polymers

There are no such residues in this entry.