



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:19 PM UTC

PDB ID : 3KJ6 / pdb_00003kj6
Title : Crystal structure of a Methylated beta2 Adrenergic Receptor-Fab complex
Authors : Bokoch, M.P.; Zou, Y.; Rasmussen, S.G.F.; Liu, C.W.; Nygaard, R.; Rosenbaum, D.M.; Fung, J.J.; Choi, H.-J.; Thian, F.S.; Kobilka, T.S.; Puglisi, J.D.; Weis, W.I.; Pardo, L.; Prosser, S.; Mueller, L.; Kobilka, B.K.
Deposited on : 2009-11-02
Resolution : 3.40 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

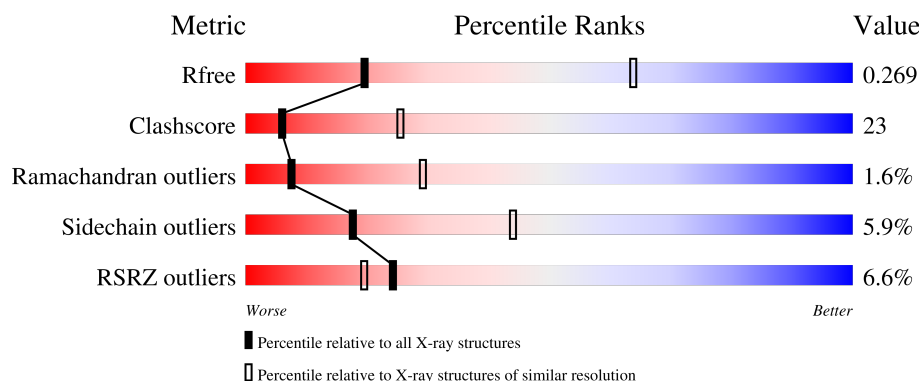
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	366	
2	L	214	
3	H	217	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4974 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 Beta-2 adrenergic receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	222	Total	C	N	O	S	0	0	0
			1656	1090	271	286	9			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP P07550
A	1	PHE	-	expression tag	UNP P07550
A	16	ARG	GLY	variant	UNP P07550
A	27	GLN	GLU	variant	UNP P07550
A	187	GLU	ASN	engineered mutation	UNP P07550

- Molecule 2 is a protein called Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	214	Total	C	N	O	S	0	0	0
			1678	1050	278	341	9			

- Molecule 3 is a protein called Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	217	Total	C	N	O	S	0	0	0
			1635	1031	269	328	7			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

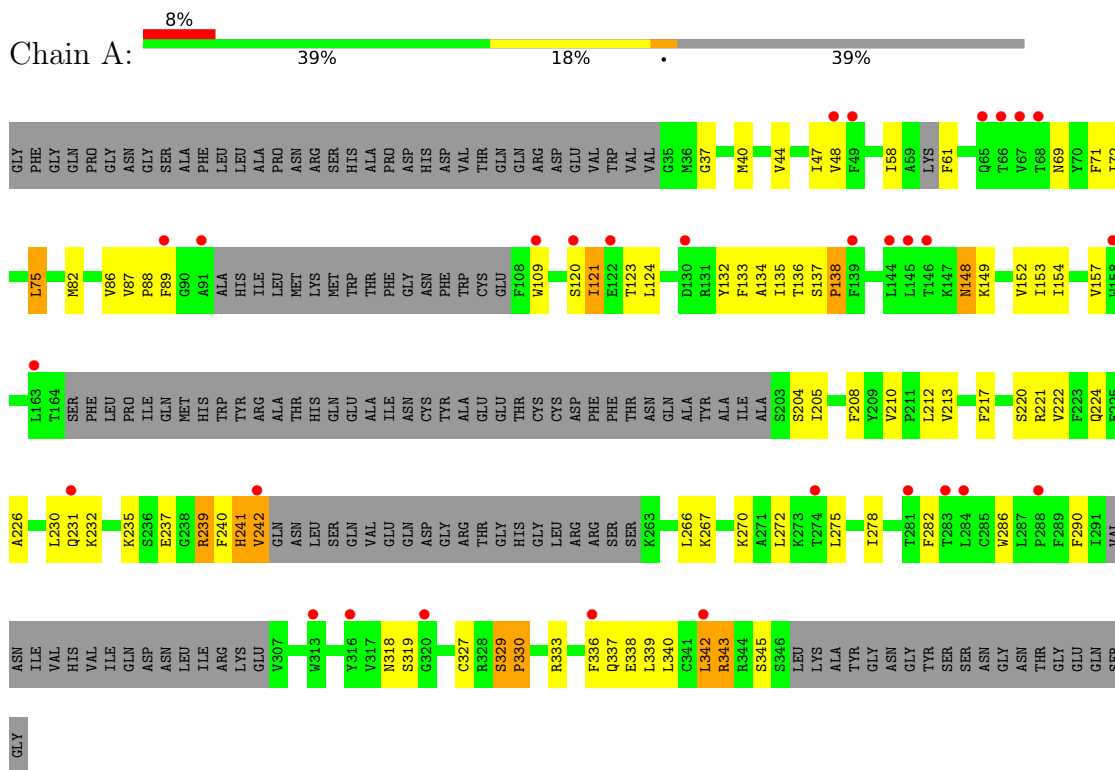


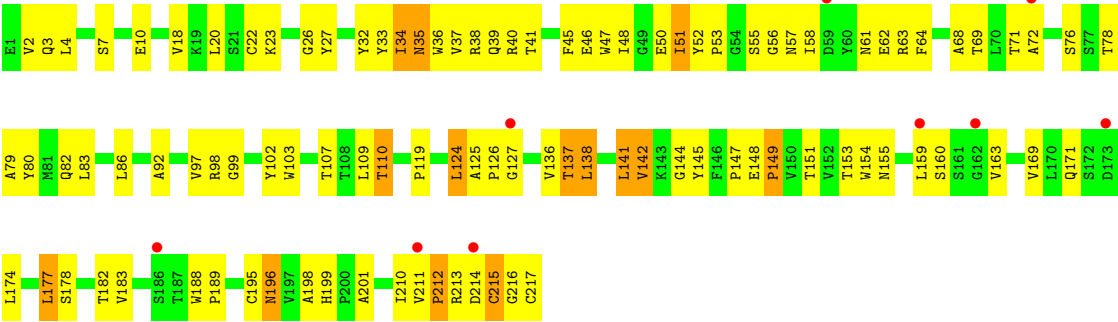
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

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: Beta-2 adrenergic receptor





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	338.60Å 48.60Å 88.90Å 90.00° 104.40° 90.00°	Depositor
Resolution (Å)	44.41 – 3.40 44.41 – 3.40	Depositor EDS
% Data completeness (in resolution range)	95.8 (44.41-3.40) 95.8 (44.41-3.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 3.40Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.238 , 0.280 0.230 , 0.269	Depositor DCC
R_{free} test set	1912 reflections (9.77%)	wwPDB-VP
Wilson B-factor (Å ²)	102.1	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 131.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.025 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4974	wwPDB-VP
Average B, all atoms (Å ²)	144.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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

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	0.29	0/1686	0.73	1/2295 (0.0%)
2	L	0.26	0/1716	0.69	0/2324
3	H	0.29	0/1677	0.76	2/2290 (0.1%)
All	All	0.28	0/5079	0.73	3/6909 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	199	HIS	CA-C-N	5.93	125.63	119.05
3	H	199	HIS	C-N-CA	5.93	125.63	119.05
1	A	232	LYS	N-CA-C	-5.49	106.56	113.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	1656	0	1631	65	0
2	L	1678	0	1610	88	0
3	H	1635	0	1578	85	0
4	A	5	0	0	0	0
All	All	4974	0	4819	222	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 23.

All (222) 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)
3:H:214:ASP:O	3:H:215:CYS:HB2	1.65	0.93
3:H:214:ASP:C	3:H:216:GLY:HA3	2.02	0.85
2:L:19:VAL:HG13	2:L:75:ILE:HG23	1.62	0.82
2:L:38:GLN:HE21	3:H:39:GLN:HE22	1.25	0.81
1:A:241:HIS:O	1:A:242:VAL:HG23	1.83	0.79
1:A:235:LYS:O	1:A:239:ARG:HD3	1.85	0.76
2:L:108:ARG:HG2	2:L:109:ALA:H	1.52	0.75
2:L:169:LYS:H	2:L:169:LYS:HD2	1.51	0.73
1:A:231:GLN:HE21	1:A:235:LYS:HD2	1.53	0.73
1:A:148:ASN:HD22	1:A:149:LYS:N	1.90	0.70
3:H:214:ASP:O	3:H:216:GLY:HA3	1.91	0.70
2:L:38:GLN:NE2	3:H:39:GLN:HE22	1.89	0.69
3:H:34:ILE:HG13	3:H:34:ILE:O	1.93	0.69
3:H:119:PRO:HB2	3:H:142:VAL:HG23	1.75	0.69
2:L:120:PRO:HD3	2:L:132:VAL:HG22	1.75	0.69
1:A:71:PHE:HD2	1:A:154:ILE:HD11	1.58	0.68
2:L:44:PRO:HG2	3:H:103:TRP:CD2	2.28	0.68
1:A:329:SER:HB2	1:A:330:PRO:HD3	1.75	0.68
3:H:69:THR:OG1	3:H:82:GLN:HB3	1.94	0.67
2:L:118:PHE:CD2	3:H:124:LEU:HB3	2.30	0.67
3:H:213:ARG:HG2	3:H:214:ASP:H	1.59	0.66
1:A:75:LEU:HD11	1:A:123:THR:HB	1.78	0.66
1:A:266:LEU:HD21	3:H:57:ASN:HB2	1.79	0.63
1:A:267:LYS:C	1:A:267:LYS:HD3	2.23	0.62
3:H:35:ASN:ND2	3:H:50:GLU:HG3	2.14	0.62
1:A:213:VAL:O	1:A:217:PHE:HD1	1.82	0.62
1:A:226:ALA:O	1:A:230:LEU:HB2	2.00	0.61
3:H:154:TRP:CD1	3:H:163:VAL:HG21	2.36	0.61
2:L:29:ILE:HG22	2:L:92:ASP:HB3	1.83	0.61
3:H:213:ARG:CG	3:H:214:ASP:H	2.14	0.61
1:A:89:PHE:HB2	1:A:109:TRP:HE1	1.66	0.60
2:L:50:ARG:HH21	2:L:53:ARG:NH1	1.98	0.60
3:H:215:CYS:N	3:H:216:GLY:HA3	2.14	0.59
2:L:144:ILE:HD13	2:L:145:ASN:H	1.66	0.59
1:A:120:SER:O	1:A:124:LEU:HD23	2.03	0.58
1:A:231:GLN:NE2	1:A:235:LYS:HD2	2.17	0.58
2:L:36:PHE:HA	2:L:47:LEU:HD22	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:PHE:HB2	1:A:109:TRP:NE1	2.19	0.57
2:L:29:ILE:HG13	2:L:71:TYR:OH	2.05	0.57
2:L:50:ARG:HH21	2:L:53:ARG:HH12	1.52	0.57
2:L:66:GLY:HA3	2:L:71:TYR:HA	1.86	0.57
2:L:138:ASN:HA	2:L:172:THR:HG21	1.86	0.57
2:L:157:ASN:C	2:L:157:ASN:HD22	2.13	0.56
3:H:110:THR:HG21	3:H:147:PRO:HB2	1.86	0.56
3:H:18:VAL:HG22	3:H:86:LEU:HD11	1.87	0.56
3:H:211:VAL:HG13	3:H:212:PRO:HD2	1.87	0.56
3:H:188:TRP:CE3	3:H:189:PRO:HD3	2.41	0.56
2:L:139:PHE:CE1	2:L:144:ILE:HB	2.41	0.56
1:A:239:ARG:HB2	3:H:99:GLY:O	2.05	0.55
1:A:69:ASN:O	1:A:72:ILE:HG22	2.07	0.55
3:H:56:GLY:O	3:H:58:ILE:HG13	2.07	0.55
2:L:108:ARG:HG2	2:L:109:ALA:N	2.22	0.55
2:L:13:ALA:O	2:L:106:ILE:HA	2.07	0.55
1:A:134:ALA:O	1:A:137:SER:HB3	2.07	0.54
2:L:137:ASN:HD22	2:L:174:SER:HB3	1.72	0.54
1:A:153:ILE:O	1:A:157:VAL:HG23	2.07	0.54
2:L:59:PRO:HG2	2:L:62:PHE:CD1	2.43	0.53
3:H:35:ASN:N	3:H:35:ASN:HD22	2.06	0.53
3:H:23:LYS:HG3	3:H:78:THR:HG22	1.91	0.53
2:L:52:ASN:HD22	2:L:53:ARG:N	2.07	0.53
3:H:52:TYR:HD2	3:H:55:SER:H	1.56	0.53
1:A:135:ILE:HD12	1:A:222:VAL:HG13	1.90	0.53
1:A:337:GLN:HG3	1:A:343:ARG:HB3	1.89	0.53
2:L:160:LEU:HD23	2:L:160:LEU:C	2.34	0.53
1:A:121:ILE:HD11	1:A:212:LEU:HB2	1.91	0.53
3:H:33:TYR:CD2	3:H:52:TYR:HB2	2.44	0.53
3:H:34:ILE:HD13	3:H:79:ALA:HB3	1.91	0.52
1:A:71:PHE:HZ	1:A:153:ILE:HD11	1.74	0.52
1:A:286:TRP:O	1:A:290:PHE:HD1	1.93	0.52
2:L:149:LYS:HG2	2:L:154:GLU:HB3	1.91	0.52
3:H:119:PRO:HB3	3:H:145:TYR:HB3	1.91	0.52
2:L:44:PRO:HG2	3:H:103:TRP:CG	2.45	0.52
2:L:193:THR:HG23	2:L:208:SER:HB2	1.92	0.52
3:H:153:THR:HG23	3:H:196:ASN:HB2	1.91	0.52
2:L:142:LYS:HD2	2:L:173:TYR:CE1	2.44	0.52
2:L:4:MET:HE3	2:L:25:ALA:HB2	1.92	0.51
2:L:6:GLN:HE22	2:L:87:TYR:HA	1.76	0.51
1:A:337:GLN:HE22	1:A:345:SER:H	1.59	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:181:LEU:HB3	2:L:185:GLU:HB2	1.92	0.51
3:H:124:LEU:HD21	3:H:141:LEU:HB2	1.92	0.51
1:A:121:ILE:HD13	1:A:121:ILE:O	2.10	0.51
2:L:138:ASN:HA	2:L:172:THR:CG2	2.40	0.51
3:H:177:LEU:HD12	3:H:178:SER:N	2.26	0.51
1:A:240:PHE:O	1:A:241:HIS:C	2.53	0.51
3:H:188:TRP:CH2	3:H:212:PRO:HA	2.46	0.51
3:H:61:ASN:HD21	3:H:63:ARG:NH2	2.10	0.50
2:L:148:TRP:O	2:L:154:GLU:HB2	2.11	0.50
1:A:235:LYS:HE3	2:L:92:ASP:O	2.11	0.50
3:H:137:THR:HG22	3:H:182:THR:HG23	1.93	0.50
3:H:142:VAL:HG13	3:H:177:LEU:HG	1.94	0.50
1:A:71:PHE:CD2	1:A:154:ILE:HD11	2.43	0.50
2:L:136:LEU:HD21	2:L:146:VAL:HG11	1.92	0.50
1:A:148:ASN:HD22	1:A:149:LYS:H	1.56	0.50
3:H:188:TRP:CZ2	3:H:212:PRO:HA	2.47	0.50
2:L:96:TYR:HD1	3:H:47:TRP:CD1	2.30	0.50
2:L:113:PRO:HG2	2:L:205:ILE:HD12	1.94	0.50
2:L:144:ILE:HD13	2:L:145:ASN:N	2.27	0.49
2:L:195:GLU:HG2	2:L:206:VAL:HG22	1.93	0.49
3:H:4:LEU:HB3	3:H:22:CYS:SG	2.52	0.49
3:H:183:VAL:HG21	3:H:188:TRP:HB2	1.95	0.49
3:H:35:ASN:HD21	3:H:50:GLU:HG3	1.78	0.49
3:H:213:ARG:HG2	3:H:214:ASP:N	2.27	0.49
2:L:155:ARG:HG2	2:L:156:GLN:H	1.77	0.49
2:L:118:PHE:CG	3:H:124:LEU:HB3	2.48	0.49
1:A:135:ILE:HD11	1:A:272:LEU:HD11	1.95	0.49
2:L:160:LEU:HD21	3:H:169:VAL:HB	1.95	0.49
1:A:275:LEU:HA	1:A:278:ILE:HG22	1.94	0.48
3:H:155:ASN:HD22	3:H:159:LEU:HD13	1.78	0.48
1:A:204:SER:O	1:A:208:PHE:HB3	2.12	0.48
1:A:337:GLN:NE2	1:A:345:SER:H	2.12	0.48
3:H:68:ALA:HA	3:H:82:GLN:O	2.12	0.48
2:L:59:PRO:HG2	2:L:62:PHE:CE1	2.47	0.48
1:A:220:SER:O	1:A:224:GLN:HG2	2.13	0.48
2:L:148:TRP:CE2	2:L:179:LEU:HD22	2.48	0.48
2:L:157:ASN:HD22	2:L:158:GLY:N	2.11	0.48
3:H:2:VAL:HA	3:H:26:GLY:HA3	1.95	0.48
3:H:38:ARG:HD2	3:H:46:GLU:OE2	2.14	0.48
3:H:51:ILE:HD13	3:H:51:ILE:C	2.39	0.48
2:L:132:VAL:HB	2:L:179:LEU:CD2	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:LYS:HB2	2:L:32:TYR:CZ	2.48	0.47
1:A:132:TYR:CE1	1:A:221:ARG:HG2	2.50	0.47
1:A:282:PHE:HD1	1:A:318:ASN:HD22	1.62	0.47
2:L:159:VAL:HG12	2:L:161:ASN:HD21	1.79	0.47
3:H:40:ARG:HA	3:H:92:ALA:HB1	1.97	0.47
3:H:136:VAL:HG12	3:H:138:LEU:HD23	1.96	0.47
2:L:24:LYS:HG2	2:L:70:ASP:HB3	1.96	0.47
1:A:87:VAL:HB	1:A:88:PRO:HD3	1.97	0.47
2:L:8:PRO:O	2:L:102:THR:HB	2.15	0.47
3:H:138:LEU:HD12	3:H:210:ILE:HD12	1.97	0.47
2:L:136:LEU:HD11	2:L:196:ALA:HB2	1.97	0.47
3:H:76:SER:O	3:H:78:THR:HG23	2.15	0.46
1:A:82:MET:O	1:A:87:VAL:HG23	2.15	0.46
1:A:132:TYR:HE1	1:A:221:ARG:HG2	1.80	0.46
2:L:2:ILE:HG23	2:L:97:THR:OG1	2.15	0.46
2:L:184:ASP:O	2:L:188:ARG:HG3	2.16	0.46
3:H:119:PRO:CB	3:H:142:VAL:HG23	2.44	0.46
2:L:16:GLY:O	2:L:77:SER:HA	2.15	0.46
2:L:198:HIS:HB3	2:L:200:THR:HG22	1.98	0.46
1:A:327:CYS:O	1:A:333:ARG:HB2	2.15	0.46
2:L:49:TYR:CD1	2:L:50:ARG:HB2	2.50	0.46
3:H:71:THR:HG23	3:H:80:TYR:HB2	1.98	0.46
2:L:36:PHE:CZ	2:L:87:TYR:HB2	2.51	0.46
1:A:47:ILE:HD12	1:A:319:SER:CB	2.46	0.46
2:L:33:LEU:HD13	2:L:71:TYR:CD1	2.51	0.46
2:L:48:ILE:HG12	2:L:54:LEU:HA	1.98	0.46
2:L:29:ILE:HG22	2:L:92:ASP:CB	2.44	0.46
3:H:151:THR:HB	3:H:198:ALA:HB3	1.98	0.46
2:L:23:CYS:HB2	2:L:35:TRP:CH2	2.51	0.46
2:L:75:ILE:O	2:L:75:ILE:HG12	2.15	0.46
2:L:108:ARG:HD2	2:L:171:SER:O	2.16	0.46
3:H:215:CYS:N	3:H:216:GLY:CA	2.80	0.45
1:A:266:LEU:HD11	3:H:57:ASN:CG	2.42	0.45
1:A:71:PHE:CZ	1:A:153:ILE:HD11	2.52	0.45
2:L:29:ILE:HG21	2:L:90:GLN:HB2	1.98	0.45
2:L:19:VAL:HG11	2:L:104:LEU:HD11	1.98	0.45
3:H:40:ARG:HA	3:H:92:ALA:CB	2.47	0.45
3:H:148:GLU:OE1	3:H:149:PRO:HA	2.17	0.45
3:H:125:ALA:HB1	3:H:126:PRO:HD2	1.99	0.45
2:L:136:LEU:HD23	2:L:175:MET:HE3	1.98	0.45
3:H:144:GLY:C	3:H:174:LEU:HD12	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:160:SER:O	3:H:163:VAL:HG12	2.17	0.44
2:L:2:ILE:HG12	2:L:2:ILE:O	2.16	0.44
2:L:29:ILE:HA	2:L:92:ASP:OD2	2.16	0.44
3:H:27:TYR:CZ	3:H:98:ARG:HD2	2.52	0.44
1:A:133:PHE:O	1:A:137:SER:HB2	2.18	0.44
3:H:38:ARG:HB2	3:H:48:ILE:HD11	1.98	0.44
2:L:115:VAL:HG11	2:L:205:ILE:HG22	2.00	0.44
3:H:34:ILE:HD11	3:H:36:TRP:HE1	1.83	0.44
1:A:47:ILE:HD12	1:A:319:SER:HB3	2.00	0.44
1:A:137:SER:OG	1:A:138:PRO:HD2	2.18	0.43
1:A:241:HIS:HB2	3:H:32:TYR:CE1	2.54	0.43
2:L:44:PRO:HG2	3:H:103:TRP:CE3	2.53	0.43
3:H:52:TYR:HA	3:H:53:PRO:HD2	1.84	0.43
2:L:115:VAL:HG13	2:L:207:LYS:HG2	2.01	0.43
1:A:132:TYR:O	1:A:136:THR:HG22	2.19	0.43
2:L:47:LEU:HA	2:L:58:VAL:HG21	2.01	0.43
2:L:149:LYS:HG2	2:L:154:GLU:CB	2.49	0.43
1:A:82:MET:HG2	1:A:86:VAL:HB	2.00	0.42
1:A:342:LEU:O	1:A:343:ARG:C	2.62	0.42
2:L:52:ASN:HD21	2:L:53:ARG:HH11	1.66	0.42
3:H:142:VAL:O	3:H:142:VAL:HG22	2.19	0.42
2:L:4:MET:HE2	2:L:4:MET:HB3	1.92	0.42
3:H:38:ARG:CZ	3:H:64:PHE:HE2	2.32	0.42
1:A:44:VAL:O	1:A:48:VAL:HG23	2.20	0.42
2:L:136:LEU:N	2:L:136:LEU:HD22	2.35	0.42
3:H:155:ASN:ND2	3:H:159:LEU:HD13	2.34	0.42
1:A:205:ILE:O	1:A:210:VAL:HG23	2.20	0.42
2:L:19:VAL:HG21	2:L:104:LEU:HD11	2.02	0.42
3:H:37:VAL:CG1	3:H:45:PHE:HB3	2.50	0.42
1:A:37:GLY:HA2	1:A:40:MET:HE2	2.02	0.42
1:A:267:LYS:HD3	1:A:267:LYS:O	2.20	0.42
1:A:154:ILE:HD13	1:A:154:ILE:HA	1.94	0.41
3:H:147:PRO:HD2	3:H:201:ALA:HB1	2.02	0.41
2:L:132:VAL:HB	2:L:179:LEU:HD21	2.02	0.41
2:L:210:ASN:HB3	2:L:213:GLU:HG2	2.02	0.41
3:H:51:ILE:HD12	3:H:72:ALA:HB2	2.01	0.41
1:A:61:PHE:HE1	1:A:339:LEU:HD21	1.86	0.41
2:L:175:MET:HG2	2:L:176:SER:N	2.35	0.41
3:H:110:THR:HG21	3:H:147:PRO:CB	2.50	0.41
3:H:83:LEU:HD12	3:H:83:LEU:HA	1.85	0.41
1:A:338:GLU:C	1:A:340:LEU:N	2.78	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:35:TRP:CD2	2:L:73:LEU:HB2	2.55	0.41
1:A:149:LYS:HA	1:A:152:VAL:HG22	2.02	0.41
2:L:37:GLN:HB2	2:L:47:LEU:HD21	2.03	0.41
2:L:137:ASN:ND2	2:L:174:SER:HB3	2.34	0.41
1:A:338:GLU:C	1:A:340:LEU:H	2.27	0.41
3:H:7:SER:HB2	3:H:107:THR:HG23	2.02	0.41
1:A:270:LYS:HD3	1:A:270:LYS:HA	1.82	0.40
1:A:336:PHE:O	1:A:340:LEU:HG	2.21	0.40
2:L:139:PHE:CD2	2:L:173:TYR:O	2.74	0.40
2:L:160:LEU:HD12	3:H:171:GLN:HB2	2.03	0.40
3:H:41:THR:O	3:H:41:THR:HG23	2.20	0.40
2:L:115:VAL:HA	2:L:135:PHE:O	2.20	0.40
3:H:10:GLU:O	3:H:109:LEU:HD12	2.21	0.40
1:A:75:LEU:HD21	1:A:124:LEU:HD22	2.03	0.40
2:L:142:LYS:HB2	2:L:173:TYR:CE2	2.56	0.40
3:H:61:ASN:O	3:H:62:GLU:C	2.65	0.40
3:H:145:TYR:CD1	3:H:145:TYR:C	2.99	0.40
2:L:159:VAL:HG12	2:L:161:ASN:ND2	2.36	0.40
1:A:148:ASN:HD22	1:A:148:ASN:C	2.28	0.40
3:H:3:GLN:HA	3:H:102:TYR:CE2	2.57	0.40
3:H:154:TRP:CZ3	3:H:195:CYS:HB3	2.56	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	210/366 (57%)	190 (90%)	15 (7%)	5 (2%)	4	22
2	L	212/214 (99%)	190 (90%)	21 (10%)	1 (0%)	24	54
3	H	215/217 (99%)	184 (86%)	27 (13%)	4 (2%)	6	26
All	All	637/797 (80%)	564 (88%)	63 (10%)	10 (2%)	7	29

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	H	215	CYS
3	H	212	PRO
1	A	329	SER
1	A	343	ARG
3	H	127	GLY
1	A	241	HIS
1	A	330	PRO
1	A	138	PRO
2	L	2	ILE
3	H	149	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	169/317 (53%)	161 (95%)	8 (5%)	23	50
2	L	191/191 (100%)	181 (95%)	10 (5%)	21	48
3	H	183/183 (100%)	169 (92%)	14 (8%)	12	37
All	All	543/691 (79%)	511 (94%)	32 (6%)	18	44

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	ILE
1	A	75	LEU
1	A	121	ILE
1	A	148	ASN
1	A	237	GLU
1	A	239	ARG
1	A	242	VAL
1	A	342	LEU
2	L	19	VAL
2	L	38	GLN
2	L	47	LEU

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Mol	Chain	Res	Type
2	L	50	ARG
2	L	52	ASN
2	L	75	ILE
2	L	102	THR
2	L	144	ILE
2	L	154	GLU
2	L	182	THR
3	H	20	LEU
3	H	34	ILE
3	H	35	ASN
3	H	51	ILE
3	H	97	VAL
3	H	110	THR
3	H	124	LEU
3	H	137	THR
3	H	138	LEU
3	H	141	LEU
3	H	142	VAL
3	H	177	LEU
3	H	196	ASN
3	H	217	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	ASN
1	A	231	GLN
1	A	312	ASN
1	A	318	ASN
1	A	337	GLN
2	L	6	GLN
2	L	30	ASN
2	L	38	GLN
2	L	52	ASN
2	L	137	ASN
2	L	157	ASN
2	L	161	ASN
2	L	198	HIS
3	H	35	ASN
3	H	133	ASN

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](#)

1 ligand is 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	$\# Z > 2$
4	SO4	A	1001	-	4,4,4	0.24	0	6,6,6	0.06	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	222/366 (60%)	0.97	30 (13%) 7 8	82, 220, 283, 334	0
2	L	214/214 (100%)	0.38	4 (1%) 66 51	65, 102, 149, 268	0
3	H	217/217 (100%)	0.42	9 (4%) 41 29	65, 104, 181, 282	0
All	All	653/797 (81%)	0.59	43 (6%) 24 19	65, 120, 264, 334	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	68	THR	4.8
1	A	139	PHE	4.2
1	A	313	TRP	4.1
1	A	67	VAL	3.8
2	L	5	THR	3.8
2	L	116	SER	3.8
1	A	242	VAL	3.5
1	A	320	GLY	3.5
1	A	109	TRP	3.4
3	H	173	ASP	3.4
1	A	274	THR	3.2
1	A	342	LEU	3.2
1	A	144	LEU	3.1
3	H	211	VAL	3.1
3	H	59	ASP	3.1
1	A	130	ASP	3.0
1	A	316	TYR	3.0
1	A	336	PHE	3.0
1	A	91	ALA	2.9
1	A	163	LEU	2.8
1	A	120	SER	2.8
1	A	146	THR	2.8
1	A	281	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	158	TRP	2.8
1	A	89	PHE	2.6
3	H	214	ASP	2.6
1	A	65	GLN	2.5
1	A	283	THR	2.4
3	H	162	GLY	2.4
3	H	186	SER	2.4
3	H	72	ALA	2.4
1	A	122	GLU	2.4
3	H	159	LEU	2.3
2	L	76	SER	2.3
1	A	231	GLN	2.3
1	A	145	LEU	2.2
1	A	49	PHE	2.2
1	A	284	LEU	2.2
3	H	127	GLY	2.2
1	A	66	THR	2.1
1	A	288	PRO	2.1
2	L	70	ASP	2.1
1	A	48	VAL	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 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($\text{\AA}^2$)	Q<0.9
4	SO4	A	1001	5/5	0.78	0.21	153,154,163,167	0

6.5 Other polymers ⓘ

There are no such residues in this entry.