



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 19, 2026 – 09:24 PM UTC

PDB ID : 4KKU / pdb_00004kku
Title : Structure of BesA (Selenomethinone derivative - P212121)
Authors : Greene, N.P.; Hinchliffe, P.; Crow, A.; Ababou, A.; Hughes, C.; Koronakis, V.
Deposited on : 2013-05-06
Resolution : 2.35 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

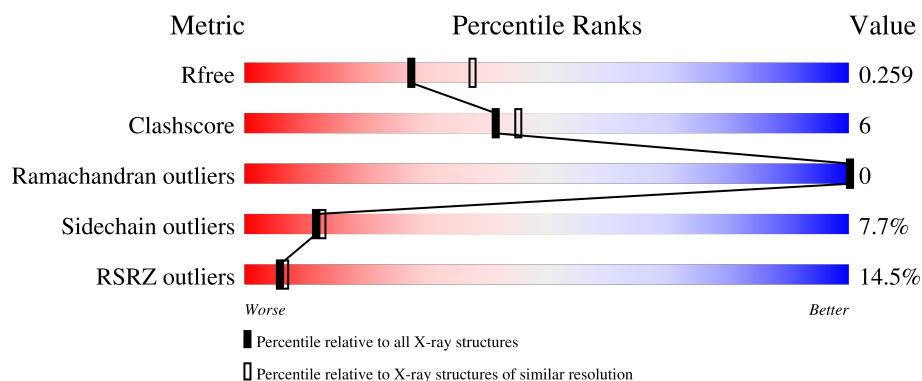
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	296	11% 74% 13% • 10%
1	B	296	8% 73% 16% • 10%
1	C	296	14% 78% 9% •• 11%
1	D	296	19% 75% 14% • 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8454 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 Membrane fusion protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	Se	0	3	0
			2075	1331	340	399	5			
1	B	267	Total	C	N	O	Se	0	3	0
			2084	1339	342	398	5			
1	C	264	Total	C	N	O	Se	0	1	0
			2043	1311	337	391	4			
1	D	266	Total	C	N	O	Se	0	3	0
			2074	1330	340	400	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	GLY	-	expression tag	UNP O51166
A	23	SER	-	expression tag	UNP O51166
A	24	HIS	-	expression tag	UNP O51166
A	25	MSE	-	expression tag	UNP O51166
B	22	GLY	-	expression tag	UNP O51166
B	23	SER	-	expression tag	UNP O51166
B	24	HIS	-	expression tag	UNP O51166
B	25	MSE	-	expression tag	UNP O51166
C	22	GLY	-	expression tag	UNP O51166
C	23	SER	-	expression tag	UNP O51166
C	24	HIS	-	expression tag	UNP O51166
C	25	MSE	-	expression tag	UNP O51166
D	22	GLY	-	expression tag	UNP O51166
D	23	SER	-	expression tag	UNP O51166
D	24	HIS	-	expression tag	UNP O51166
D	25	MSE	-	expression tag	UNP O51166

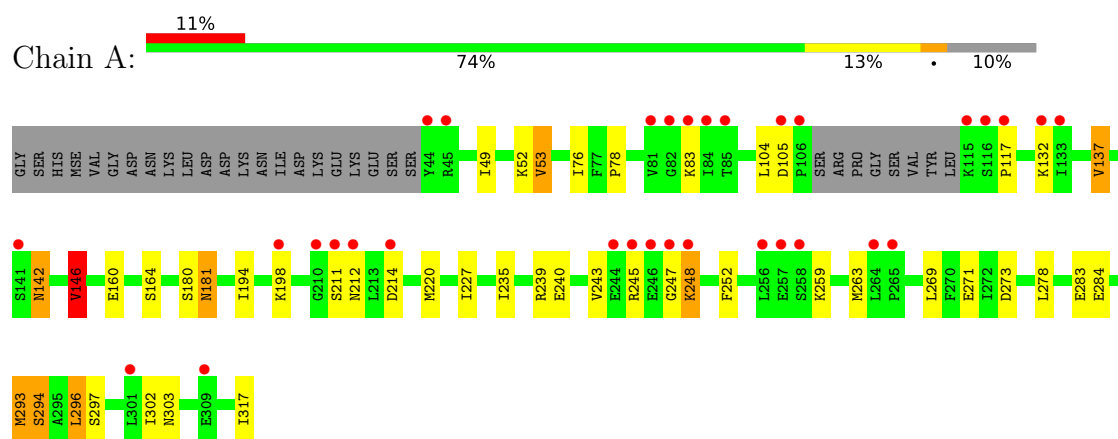
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	50	Total 50	O 50	0	0
2	B	47	Total 47	O 47	0	0
2	C	37	Total 37	O 37	0	0
2	D	44	Total 44	O 44	0	0

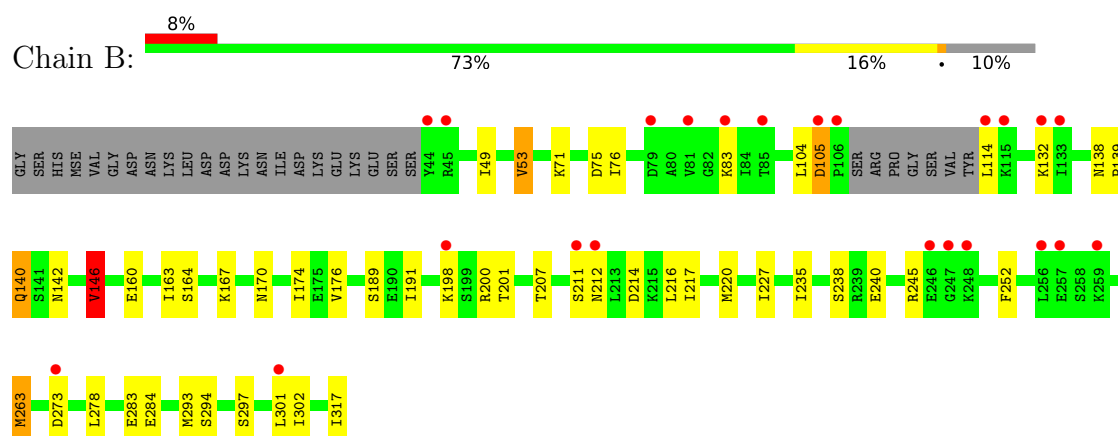
3 Residue-property plots [i](#)

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.

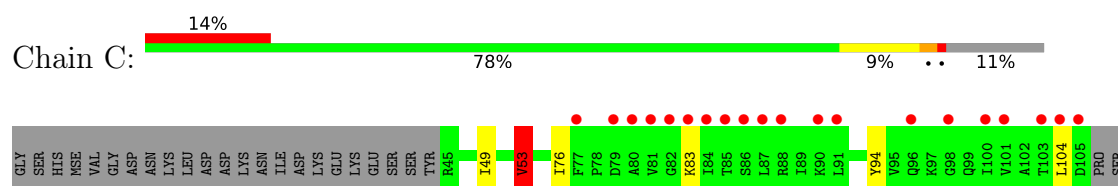
- Molecule 1: Membrane fusion protein

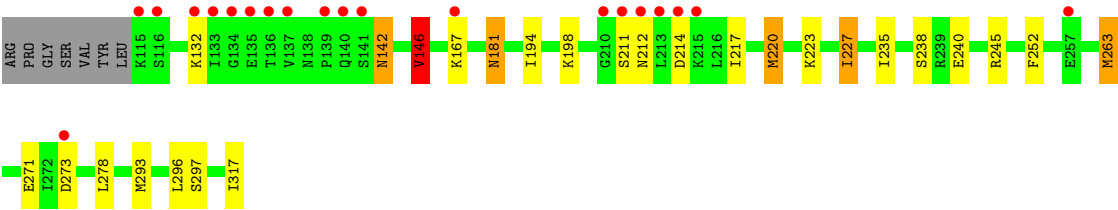


- Molecule 1: Membrane fusion protein

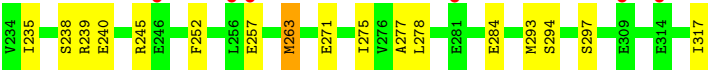
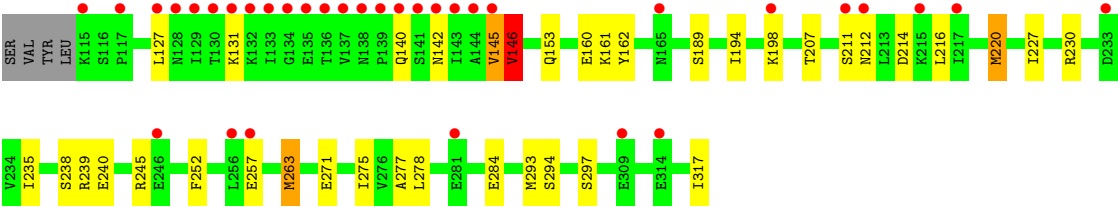
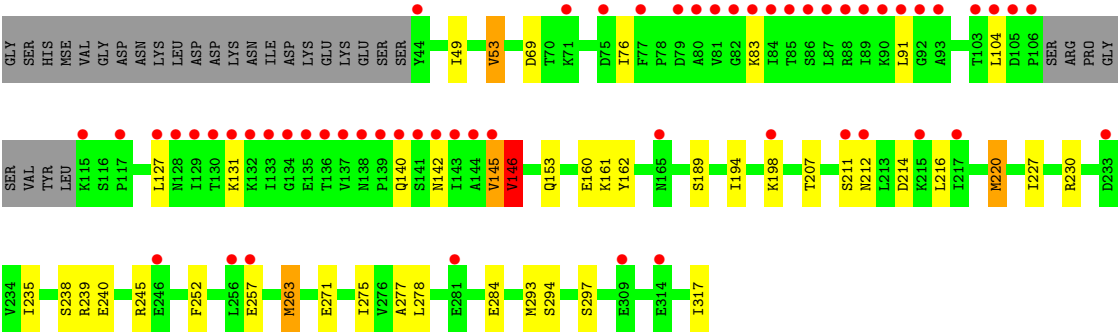


- Molecule 1: Membrane fusion protein





● Molecule 1: Membrane fusion protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.74Å 152.13Å 155.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	77.74 – 2.35 77.74 – 2.35	Depositor EDS
% Data completeness (in resolution range)	98.5 (77.74-2.35) 98.5 (77.74-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.236 , 0.264 (Not available) , 0.259	Depositor DCC
R_{free} test set	3697 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 23.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8454	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 54.73 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.5035e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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](#)

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	1.22	2/2094 (0.1%)	1.07	3/2814 (0.1%)
1	B	1.25	5/2103 (0.2%)	1.09	6/2825 (0.2%)
1	C	1.16	1/2060 (0.0%)	1.09	4/2766 (0.1%)
1	D	1.11	3/2093 (0.1%)	1.10	7/2813 (0.2%)
All	All	1.19	11/8350 (0.1%)	1.09	20/11218 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	191	ILE	CA-CB	-7.54	1.44	1.54
1	B	227	ILE	CA-CB	-6.44	1.46	1.54
1	B	71	LYS	N-CA	6.03	1.53	1.46
1	C	227	ILE	C-O	-5.90	1.18	1.24
1	D	277	ALA	CA-C	-5.57	1.46	1.52
1	B	200	ARG	CA-C	-5.57	1.45	1.53
1	D	275	ILE	C-O	-5.51	1.18	1.24
1	D	162	TYR	C-O	-5.29	1.17	1.24
1	B	201	THR	CA-C	-5.23	1.46	1.52
1	A	180	SER	C-O	-5.22	1.17	1.24
1	A	303	ASN	CA-C	5.21	1.59	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	220	MSE	CA-CB-CG	11.24	136.59	114.10
1	D	220	MSE	N-CA-CB	-10.63	94.13	110.29
1	D	220	MSE	CG-SE-CE	-9.68	77.64	98.92
1	A	146	VAL	CB-CA-C	-9.05	97.05	110.81
1	B	146	VAL	CB-CA-C	-8.94	97.23	110.81
1	C	146	VAL	CB-CA-C	-8.41	98.02	110.81
1	D	146	VAL	CB-CA-C	-7.83	98.91	110.81
1	B	105	ASP	CB-CA-C	7.37	120.00	110.13
1	C	263	MSE	CG-SE-CE	5.94	111.99	98.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	140	GLN	CA-CB-CG	5.91	125.91	114.10
1	D	263	MSE	CG-SE-CE	5.74	111.55	98.92
1	D	194	ILE	CB-CA-C	-5.68	103.87	110.91
1	A	194	ILE	CB-CA-C	-5.63	103.93	110.91
1	B	263	MSE	CG-SE-CE	5.39	110.78	98.92
1	C	194	ILE	CB-CA-C	-5.37	104.25	110.91
1	B	174	ILE	CB-CA-C	-5.26	103.77	110.98
1	A	105	ASP	CB-CA-C	-5.23	104.14	110.15
1	C	53	VAL	N-CA-CB	-5.18	103.16	110.26
1	D	220	MSE	CB-CA-C	5.08	118.81	109.46
1	B	163	ILE	N-CA-C	5.07	117.43	111.09

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	2075	0	2199	36	0
1	B	2084	0	2219	33	0
1	C	2043	0	2180	24	0
1	D	2074	0	2197	26	0
2	A	50	0	0	7	0
2	B	47	0	0	4	0
2	C	37	0	0	6	0
2	D	44	0	0	5	0
All	All	8454	0	8795	109	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 6.

All (109) 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:D:230:ARG:HG3	2:D:420:HOH:O	1.49	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:245:ARG:NH2	1:D:263:MSE:HE1	1.91	0.86
1:A:142:ASN:ND2	2:A:412:HOH:O	2.09	0.84
1:B:245:ARG:NH2	1:B:263:MSE:HE1	1.93	0.83
1:D:69:ASP:OD2	2:D:441:HOH:O	1.99	0.80
1:C:142:ASN:ND2	2:C:416:HOH:O	2.14	0.80
1:C:252:PHE:HZ	1:C:263:MSE:HE2	1.48	0.77
1:A:220[A]:MSE:HE1	1:B:220[A]:MSE:CE	2.15	0.75
1:C:252:PHE:HZ	1:C:263:MSE:CE	1.98	0.75
1:A:293:MSE:HA	1:A:296:LEU:HD22	1.70	0.74
1:B:252:PHE:HZ	1:B:263:MSE:HE2	1.54	0.73
1:C:252:PHE:CZ	1:C:263:MSE:CE	2.71	0.73
1:D:252:PHE:HZ	1:D:263:MSE:HE2	1.53	0.72
1:C:252:PHE:CZ	1:C:263:MSE:HE2	2.26	0.71
1:D:252:PHE:HZ	1:D:263:MSE:CE	2.04	0.69
1:D:252:PHE:CZ	1:D:263:MSE:CE	2.75	0.69
1:B:273:ASP:OD2	2:B:447:HOH:O	2.11	0.69
1:B:252:PHE:HZ	1:B:263:MSE:CE	2.06	0.68
1:D:252:PHE:CZ	1:D:263:MSE:HE2	2.28	0.68
1:B:252:PHE:CZ	1:B:263:MSE:HE2	2.28	0.68
1:B:301:LEU:HG	2:B:439:HOH:O	1.94	0.68
1:A:78:PRO:HG3	1:A:137:VAL:HG11	1.74	0.67
1:D:239:ARG:NH2	2:D:437:HOH:O	2.14	0.67
1:A:220[A]:MSE:CE	1:B:220[A]:MSE:CE	2.71	0.67
1:A:245:ARG:NH2	1:A:296:LEU:O	2.28	0.66
1:B:252:PHE:CZ	1:B:263:MSE:CE	2.78	0.66
1:C:223:LYS:HE3	2:C:419:HOH:O	1.95	0.66
1:A:220[A]:MSE:HE1	1:B:220[A]:MSE:HE2	1.77	0.66
1:A:273:ASP:OD2	2:A:406:HOH:O	2.15	0.65
1:A:252:PHE:HZ	1:A:263:MSE:HE2	1.61	0.65
1:A:252:PHE:HZ	1:A:263:MSE:CE	2.10	0.64
1:A:252:PHE:CZ	1:A:263:MSE:CE	2.81	0.64
1:C:245:ARG:NH1	2:C:430:HOH:O	2.31	0.64
1:A:294:SER:OG	2:A:434:HOH:O	2.12	0.64
1:A:220[A]:MSE:HE1	1:B:176:VAL:HB	1.80	0.63
1:C:245:ARG:NH2	1:C:296:LEU:O	2.32	0.62
1:A:220[A]:MSE:CE	1:B:220[A]:MSE:HE1	2.29	0.62
1:A:252:PHE:CZ	1:A:263:MSE:HE2	2.34	0.62
1:D:127:LEU:HB2	1:D:145:VAL:HG22	1.81	0.62
1:D:153:GLN:NE2	2:D:441:HOH:O	2.32	0.62
1:A:220[A]:MSE:HE1	1:B:220[A]:MSE:HE1	1.82	0.60
1:D:127:LEU:HB2	1:D:145:VAL:CG2	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:PRO:CG	1:A:137:VAL:CG1	2.80	0.60
1:C:252:PHE:CZ	1:C:263:MSE:HE3	2.36	0.59
1:A:78:PRO:HG2	1:A:137:VAL:CG1	2.34	0.57
1:D:245:ARG:NH2	1:D:263:MSE:CE	2.64	0.57
1:B:245:ARG:NH2	1:B:263:MSE:CE	2.67	0.57
1:D:189[B]:SER:HB3	1:D:207:THR:HG1	1.69	0.56
2:A:407:HOH:O	1:D:161:LYS:HD3	2.04	0.56
1:B:76:ILE:HD11	1:B:146:VAL:HG22	1.86	0.56
1:A:76:ILE:HD11	1:A:146:VAL:HG22	1.88	0.56
1:A:160:GLU:OE2	1:D:271:GLU:OE1	2.25	0.55
1:D:252:PHE:CZ	1:D:263:MSE:HE3	2.41	0.55
1:D:76:ILE:HD11	1:D:146:VAL:HG22	1.87	0.55
1:B:212:ASN:N	2:B:429:HOH:O	2.39	0.53
1:C:181:ASN:H	1:C:181:ASN:HD22	1.57	0.53
1:D:220:MSE:HE1	2:D:405:HOH:O	2.08	0.52
1:A:78:PRO:HG3	1:A:137:VAL:CG1	2.38	0.52
1:B:252:PHE:CZ	1:B:263:MSE:HE3	2.45	0.52
1:C:76:ILE:HD11	1:C:146:VAL:HG22	1.91	0.52
1:A:252:PHE:CZ	1:A:263:MSE:HE3	2.45	0.50
1:D:278:LEU:C	1:D:278:LEU:HD12	2.35	0.50
1:B:217:ILE:HD12	1:B:220[B]:MSE:HE3	1.93	0.50
1:A:117:PRO:HD2	2:A:410:HOH:O	2.11	0.50
1:C:53:VAL:HG13	1:C:235:ILE:HD12	1.94	0.49
1:A:181:ASN:HD22	1:A:181:ASN:H	1.59	0.49
1:C:217:ILE:O	1:C:220:MSE:HB2	2.13	0.49
1:C:94:TYR:OH	2:C:413:HOH:O	2.15	0.48
1:D:53:VAL:HG13	1:D:235:ILE:HD12	1.94	0.48
1:B:278:LEU:HD12	1:B:278:LEU:C	2.39	0.48
1:A:278:LEU:C	1:A:278:LEU:HD12	2.37	0.47
1:C:278:LEU:HD12	1:C:278:LEU:C	2.40	0.47
1:A:78:PRO:HG2	1:A:137:VAL:HG13	1.95	0.47
1:B:75:ASP:CG	1:B:142:ASN:ND2	2.74	0.46
1:C:223:LYS:CE	2:C:419:HOH:O	2.56	0.46
1:A:245:ARG:HE	1:A:263:MSE:HE1	1.80	0.46
1:B:75:ASP:CG	1:B:142:ASN:HD21	2.24	0.45
1:B:142:ASN:OD1	2:B:416:HOH:O	2.20	0.45
1:D:83:LYS:O	1:D:104:LEU:HA	2.17	0.44
1:A:53:VAL:HG13	1:A:235:ILE:HD12	1.99	0.44
1:D:216:LEU:HD12	1:D:216:LEU:HA	1.88	0.44
1:B:138:ASN:HB2	1:B:139:PRO:CD	2.48	0.44
1:A:239:ARG:NH2	2:A:433:HOH:O	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:GLY:O	1:A:248:LYS:HD2	2.18	0.43
1:B:167:LYS:H	1:B:170:ASN:ND2	2.16	0.43
1:B:53:VAL:HG13	1:B:235:ILE:HD12	1.99	0.43
1:C:252:PHE:CE1	1:C:263:MSE:HE3	2.53	0.43
1:A:83:LYS:O	1:A:104:LEU:HA	2.18	0.43
1:D:252:PHE:CE1	1:D:263:MSE:HE3	2.54	0.43
1:C:83:LYS:O	1:C:104:LEU:HA	2.19	0.43
2:A:407:HOH:O	1:D:161:LYS:CD	2.65	0.42
1:A:302:ILE:HD12	1:A:302:ILE:C	2.44	0.42
1:B:211:SER:O	1:B:212:ASN:CB	2.68	0.42
1:C:167[B]:LYS:HB2	1:C:167[B]:LYS:HE3	1.84	0.42
1:A:271:GLU:OE2	1:D:160:GLU:OE2	2.38	0.42
1:C:273:ASP:OD2	2:C:435:HOH:O	2.22	0.42
1:B:160:GLU:OE2	1:C:271:GLU:OE1	2.38	0.41
1:B:283:GLU:O	1:B:284:GLU:C	2.63	0.41
1:A:283:GLU:O	1:A:284:GLU:C	2.63	0.41
1:B:189[B]:SER:HB3	1:B:207:THR:HG1	1.84	0.41
1:C:181:ASN:HD22	1:C:181:ASN:N	2.18	0.41
1:A:211:SER:O	1:A:212:ASN:CB	2.67	0.41
1:B:83:LYS:O	1:B:104:LEU:HA	2.19	0.41
1:D:211:SER:O	1:D:212:ASN:CB	2.69	0.41
1:A:220[A]:MSE:HE2	1:B:220[A]:MSE:CE	2.50	0.41
1:B:302:ILE:C	1:B:302:ILE:HD12	2.46	0.41
1:C:211:SER:O	1:C:212:ASN:CB	2.68	0.41
1:C:245:ARG:HE	1:C:263:MSE:HE1	1.86	0.40
1:B:216:LEU:HD12	1:B:216:LEU:HA	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	70/296 (24%)	65 (93%)	5 (7%)	0	100	100
1	B	266/296 (90%)	257 (97%)	9 (3%)	0	100	100
1	C	261/296 (88%)	254 (97%)	7 (3%)	0	100	100
1	D	265/296 (90%)	257 (97%)	8 (3%)	0	100	100
All	All	862/1184 (73%)	833 (97%)	29 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	238/257 (93%)	215 (90%)	23 (10%)	8	7
1	B	239/257 (93%)	223 (93%)	16 (7%)	15	17
1	C	234/257 (91%)	219 (94%)	15 (6%)	16	18
1	D	238/257 (93%)	219 (92%)	19 (8%)	11	12
All	All	949/1028 (92%)	876 (92%)	73 (8%)	12	13

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ILE
1	A	52	LYS
1	A	53	VAL
1	A	132	LYS
1	A	137	VAL
1	A	142	ASN
1	A	146	VAL
1	A	164	SER
1	A	181	ASN
1	A	198	LYS
1	A	214[A]	ASP
1	A	214[B]	ASP
1	A	227	ILE

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Mol	Chain	Res	Type
1	A	240	GLU
1	A	243	VAL
1	A	248	LYS
1	A	259	LYS
1	A	269	LEU
1	A	293	MSE
1	A	294	SER
1	A	296	LEU
1	A	297	SER
1	A	317	ILE
1	B	49	ILE
1	B	53	VAL
1	B	105	ASP
1	B	114	LEU
1	B	132	LYS
1	B	140	GLN
1	B	146	VAL
1	B	164	SER
1	B	198	LYS
1	B	214	ASP
1	B	238	SER
1	B	240	GLU
1	B	293	MSE
1	B	294	SER
1	B	297	SER
1	B	317	ILE
1	C	49	ILE
1	C	53	VAL
1	C	132	LYS
1	C	142	ASN
1	C	146	VAL
1	C	181	ASN
1	C	198	LYS
1	C	214	ASP
1	C	220	MSE
1	C	227	ILE
1	C	238	SER
1	C	240	GLU
1	C	293	MSE
1	C	297	SER
1	C	317	ILE
1	D	49	ILE

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Mol	Chain	Res	Type
1	D	53	VAL
1	D	91	LEU
1	D	131	LYS
1	D	140	GLN
1	D	142	ASN
1	D	145	VAL
1	D	146	VAL
1	D	198	LYS
1	D	214	ASP
1	D	227	ILE
1	D	238	SER
1	D	240	GLU
1	D	257	GLU
1	D	284	GLU
1	D	293	MSE
1	D	294	SER
1	D	297	SER
1	D	317	ILE

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

Mol	Chain	Res	Type
1	A	181	ASN
1	B	96	GLN
1	B	142	ASN
1	B	165	ASN
1	B	170	ASN
1	C	181	ASN
1	C	274	ASN
1	C	285	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	262/296 (88%)	0.60	32 (12%) 8 10	11, 35, 72, 115	2 (0%)
1	B	263/296 (88%)	0.50	23 (8%) 16 18	12, 33, 68, 98	2 (0%)
1	C	260/296 (87%)	0.72	40 (15%) 5 6	18, 37, 96, 112	1 (0%)
1	D	262/296 (88%)	1.06	57 (21%) 2 2	18, 39, 107, 130	3 (1%)
All	All	1047/1184 (88%)	0.72	152 (14%) 6 7	11, 36, 90, 130	8 (0%)

All (152) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	114	LEU	8.2
1	D	106	PRO	8.1
1	A	106	PRO	7.1
1	D	91	LEU	6.9
1	D	81	VAL	6.9
1	D	130	THR	6.2
1	D	77	PHE	6.1
1	D	133	ILE	6.1
1	D	132	LYS	5.7
1	D	129	ILE	5.7
1	D	44	TYR	5.6
1	C	115	LYS	5.5
1	A	44	TYR	5.4
1	A	115	LYS	5.3
1	B	44	TYR	5.3
1	B	115	LYS	5.0
1	D	136	THR	4.9
1	C	136	THR	4.9
1	C	81	VAL	4.8
1	D	84	ILE	4.8
1	D	212	ASN	4.7

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Mol	Chain	Res	Type	RSRZ
1	C	133	ILE	4.7
1	A	211	SER	4.4
1	A	212	ASN	4.4
1	C	90	LYS	4.3
1	C	212	ASN	4.3
1	D	80	ALA	4.3
1	A	45	ARG	4.2
1	C	85	THR	4.2
1	D	139	PRO	4.1
1	B	259[A]	LYS	4.1
1	C	105	ASP	4.1
1	D	92	GLY	4.0
1	A	265	PRO	4.0
1	D	134	GLY	4.0
1	C	80	ALA	3.9
1	D	215	LYS	3.9
1	B	248	LYS	3.9
1	D	142	ASN	3.9
1	B	106	PRO	3.8
1	D	104	LEU	3.8
1	D	141	SER	3.8
1	D	83	LYS	3.7
1	B	132	LYS	3.7
1	D	127	LEU	3.6
1	A	248	LYS	3.6
1	C	140	GLN	3.6
1	D	128	ASN	3.6
1	C	132	LYS	3.6
1	C	211	SER	3.6
1	D	131	LYS	3.6
1	D	105	ASP	3.5
1	C	101	VAL	3.5
1	B	45	ARG	3.5
1	B	246	GLU	3.4
1	A	82	GLY	3.4
1	D	71	LYS	3.4
1	A	116	SER	3.4
1	B	247	GLY	3.4
1	C	84	ILE	3.3
1	C	86	SER	3.3
1	C	83	LYS	3.2
1	A	105	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	85	THR	3.2
1	A	264	LEU	3.1
1	D	144	ALA	3.1
1	A	83	LYS	3.1
1	B	212	ASN	3.1
1	C	215	LYS	3.0
1	D	90	LYS	3.0
1	D	93	ALA	3.0
1	C	137	VAL	3.0
1	D	89	ILE	3.0
1	D	143	ILE	3.0
1	A	245	ARG	3.0
1	D	115	LYS	3.0
1	A	117	PRO	3.0
1	D	137	VAL	2.9
1	C	98	GLY	2.9
1	C	100	ILE	2.9
1	D	82	GLY	2.9
1	A	244	GLU	2.9
1	A	85	THR	2.9
1	A	198	LYS	2.8
1	D	87	LEU	2.8
1	B	83	LYS	2.8
1	C	79	ASP	2.8
1	D	135	GLU	2.8
1	B	273	ASP	2.8
1	A	301	LEU	2.8
1	D	75	ASP	2.7
1	D	103	THR	2.7
1	D	140	GLN	2.7
1	D	257	GLU	2.7
1	B	79	ASP	2.7
1	A	246	GLU	2.7
1	B	85	THR	2.7
1	A	247	GLY	2.7
1	C	134	GLY	2.7
1	C	87	LEU	2.7
1	C	104	LEU	2.7
1	C	77	PHE	2.6
1	A	214[A]	ASP	2.6
1	D	145	VAL	2.6
1	D	281	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	309[A]	GLU	2.6
1	A	133	ILE	2.6
1	B	211	SER	2.6
1	C	139	PRO	2.6
1	C	141	SER	2.6
1	B	301	LEU	2.6
1	D	233	ASP	2.5
1	C	103	THR	2.5
1	D	86	SER	2.5
1	D	211	SER	2.5
1	C	135	GLU	2.5
1	D	79	ASP	2.5
1	A	210	GLY	2.5
1	D	138	ASN	2.4
1	C	214	ASP	2.4
1	C	257	GLU	2.4
1	A	81	VAL	2.4
1	A	257	GLU	2.4
1	D	198	LYS	2.3
1	B	257	GLU	2.3
1	D	256	LEU	2.3
1	D	117	PRO	2.3
1	B	198	LYS	2.3
1	C	210	GLY	2.3
1	A	309	GLU	2.2
1	C	116	SER	2.2
1	A	84	ILE	2.2
1	D	165	ASN	2.2
1	C	82	GLY	2.2
1	D	314	GLU	2.2
1	D	217	ILE	2.1
1	C	167[A]	LYS	2.1
1	A	256	LEU	2.1
1	C	91	LEU	2.1
1	D	246	GLU	2.1
1	C	96	GLN	2.1
1	C	88	ARG	2.1
1	A	141	SER	2.1
1	B	256	LEU	2.1
1	B	81	VAL	2.1
1	A	258	SER	2.1
1	D	88	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	213	LEU	2.0
1	A	132	LYS	2.0
1	B	105	ASP	2.0
1	C	273	ASP	2.0
1	B	133	ILE	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.