



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:31 AM UTC

PDB ID : 5V1D / pdb_00005v1d
Title : Complex structure of the bovine PERK luminal domain and its substrate peptide
Authors : Wang, P.; Li, J.; Sha, B.
Deposited on : 2017-03-02
Resolution : 2.80 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
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

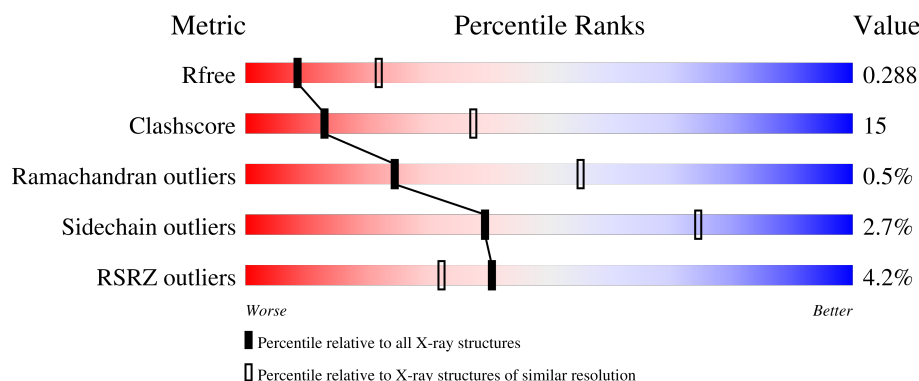
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	326	 4% 39% 19% 41%
1	B	326	 3% 36% 21% 42%
1	C	326	 2% 52% 22% 25%
1	D	326	 2% 53% 22% 24%
2	E	12	 58% 17% 25%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	F	12	17%17%50%8%25%
2	G	12	17%42%25%8%25%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7270 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 eIF2AK3 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1511	975	254	277	5			
1	B	190	Total	C	N	O	S	0	0	0
			1477	949	251	272	5			
1	C	244	Total	C	N	O	S	0	0	0
			1916	1219	321	368	8			
1	D	247	Total	C	N	O	S	0	0	0
			1944	1237	328	371	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	337	SER	CYS	engineered mutation	UNP A5D791
B	337	SER	CYS	engineered mutation	UNP A5D791
C	337	SER	CYS	engineered mutation	UNP A5D791
D	337	SER	CYS	engineered mutation	UNP A5D791

- Molecule 2 is a protein called 12-mer peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	9	Total	C	N	O	0	0	0
			83	58	14	11			
2	F	9	Total	C	N	O	0	0	0
			83	58	14	11			
2	G	9	Total	C	N	O	0	0	0
			83	58	14	11			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	43	Total	O	0	0
			43	43		

Continued on next page...

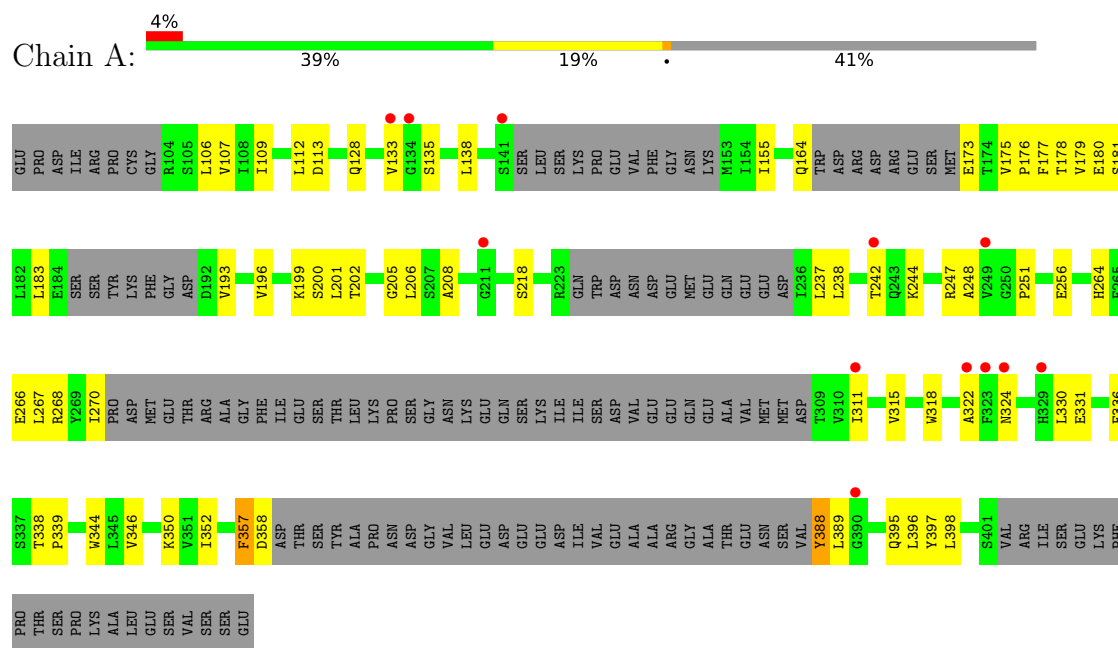
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	2	Total 2	O 2	0	0
3	B	24	Total 24	O 24	0	0
3	C	53	Total 53	O 53	0	0
3	F	1	Total 1	O 1	0	0
3	D	46	Total 46	O 46	0	0
3	G	4	Total 4	O 4	0	0

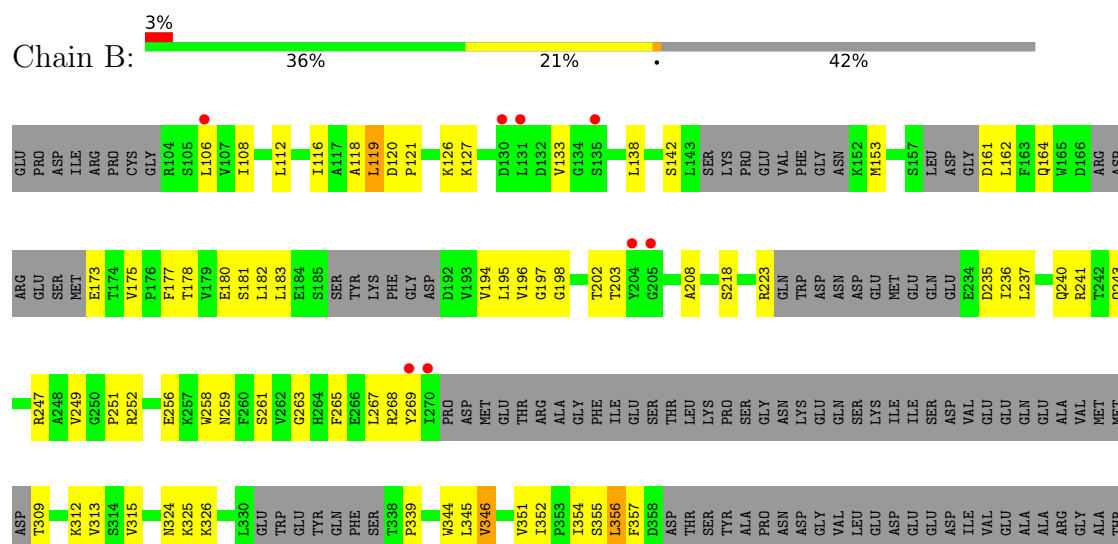
3 Residue-property plots

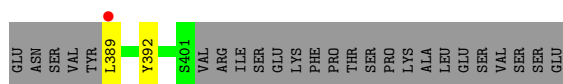
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: eIF2AK3 protein

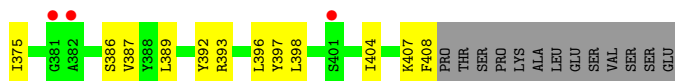
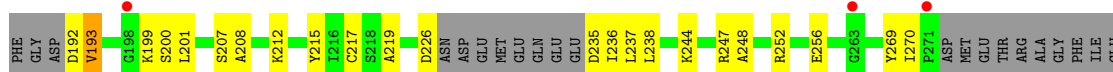
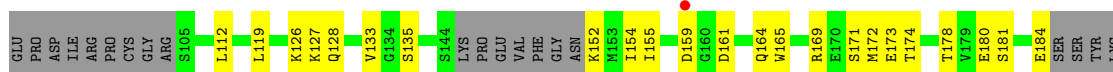


• Molecule 1: eIF2AK3 protein





• Molecule 1: eIF2AK3 protein



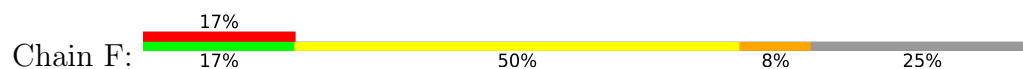
• Molecule 1: eIF2AK3 protein



• Molecule 2: 12-mer peptide



• Molecule 2: 12-mer peptide





● Molecule 2: 12-mer peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.75Å 161.55Å 104.78Å 90.00° 106.84° 90.00°	Depositor
Resolution (Å)	47.89 – 2.80 47.89 – 2.80	Depositor EDS
% Data completeness (in resolution range)	91.4 (47.89-2.80) 91.4 (47.89-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	124.80 (at 2.81Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.240 , 0.291 0.241 , 0.288	Depositor DCC
R_{free} test set	2024 reflections (4.39%)	wwPDB-VP
Wilson B-factor (Å ²)	59.7	Xtriage
Anisotropy	0.270	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 82.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.419 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	7270	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.97% 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](#)

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	0.22	0/1538	0.41	0/2073
1	B	0.20	0/1499	0.41	0/2016
1	C	0.22	0/1950	0.38	0/2632
1	D	0.29	0/1979	0.44	0/2670
2	E	0.35	0/89	0.58	0/122
2	F	0.26	0/89	0.60	0/122
2	G	0.32	0/89	0.54	0/122
All	All	0.24	0/7233	0.42	0/9757

There are no bond length outliers.

There are no bond angle outliers.

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	1511	0	1528	50	0
1	B	1477	0	1504	53	0
1	C	1916	0	1899	54	0
1	D	1944	0	1934	59	0
2	E	83	0	75	4	0
2	F	83	0	75	8	0
2	G	83	0	75	4	0
3	A	43	0	0	6	0
3	B	24	0	0	5	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	53	0	0	4	0
3	D	46	0	0	10	0
3	E	2	0	0	1	0
3	F	1	0	0	0	0
3	G	4	0	0	0	0
All	All	7270	0	7090	213	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 15.

All (213) 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:C:389:LEU:HD23	2:F:5:PRO:HA	1.62	0.81
1:A:112:LEU:HD12	1:A:339:PRO:HB2	1.65	0.79
1:D:266:GLU:HB3	1:D:268:ARG:HH22	1.49	0.77
1:D:359:ASP:HA	1:D:404:ILE:HG23	1.66	0.76
1:C:386:SER:HB3	2:F:8:PHE:HB3	1.68	0.74
1:C:165:TRP:HB2	1:C:172:MET:HE3	1.70	0.72
1:C:350:LYS:NZ	3:C:501:HOH:O	2.22	0.72
1:D:152:LYS:NZ	3:D:502:HOH:O	2.22	0.72
1:D:408:PHE:HD1	3:D:507:HOH:O	1.72	0.71
1:D:357:PHE:HB3	1:D:404:ILE:HG21	1.72	0.70
1:A:202:THR:HG23	1:A:218:SER:HB3	1.73	0.69
1:C:155:ILE:HD12	1:C:396:LEU:HB2	1.74	0.69
1:B:178:THR:H	1:B:181:SER:HB2	1.57	0.68
1:C:236:ILE:HD11	1:C:270:ILE:HD12	1.73	0.68
1:A:178:THR:H	1:A:181:SER:HB2	1.59	0.67
1:B:202:THR:HG23	1:B:218:SER:HB3	1.78	0.66
1:D:192:ASP:N	3:D:505:HOH:O	2.27	0.66
1:D:377:GLU:HG2	1:D:404:ILE:HD11	1.77	0.66
1:B:119:LEU:HD22	1:B:127:LYS:HA	1.79	0.64
1:C:192:ASP:N	3:C:504:HOH:O	2.31	0.64
1:C:208:ALA:HA	1:C:237:LEU:HG	1.80	0.63
2:F:6:TRP:O	2:F:7:ARG:NH1	2.33	0.62
1:B:112:LEU:HD22	1:B:339:PRO:HB2	1.81	0.62
1:B:197:GLY:HA2	1:C:219:ALA:HB2	1.80	0.62
1:B:249:VAL:HG12	1:B:256:GLU:HA	1.81	0.62
1:B:252:ARG:HH21	1:C:235:ASP:N	1.99	0.60
1:A:389:LEU:HD11	1:A:396:LEU:HB3	1.83	0.60
1:A:196:VAL:HG23	1:A:247:ARG:HB2	1.84	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:LEU:HB3	1:C:201:LEU:HD13	1.83	0.60
1:D:104:ARG:NE	1:D:104:ARG:O	2.35	0.59
1:B:108:ILE:HG23	1:B:116:ILE:HD11	1.83	0.59
1:B:256:GLU:OE1	1:B:259:ASN:ND2	2.36	0.59
1:C:392:TYR:OH	1:C:393:ARG:NH1	2.35	0.59
1:B:344:TRP:NE1	3:B:504:HOH:O	2.36	0.59
1:C:319:LYS:HG2	1:C:335:GLN:HA	1.85	0.59
1:D:113:ASP:H	1:D:338:THR:HG21	1.68	0.59
1:D:375:ILE:HA	1:D:408:PHE:HE1	1.68	0.58
1:A:205:GLY:N	3:A:510:HOH:O	2.36	0.58
1:B:345:LEU:N	1:B:352:ILE:O	2.30	0.58
1:D:109:ILE:HD13	1:D:320:VAL:HG21	1.84	0.58
1:B:196:VAL:HG21	1:C:217:CYS:HB3	1.86	0.58
1:B:175:VAL:HG12	1:B:177:PHE:H	1.68	0.58
1:C:159:ASP:O	1:C:244:LYS:NZ	2.37	0.58
1:B:138:LEU:O	1:B:263:GLY:N	2.35	0.57
1:D:361:SER:HB2	1:D:406:GLU:H	1.69	0.57
1:A:176:PRO:HB2	3:A:523:HOH:O	2.03	0.57
1:B:309:THR:N	3:B:505:HOH:O	2.37	0.57
1:B:153:MET:SD	1:B:153:MET:N	2.76	0.57
1:A:133:VAL:HG23	1:A:135:SER:H	1.68	0.56
1:C:375:ILE:HA	1:C:408:PHE:HB3	1.87	0.56
1:C:126:LYS:NZ	1:C:128:GLN:OE1	2.39	0.55
1:B:389:LEU:N	3:B:506:HOH:O	2.39	0.54
1:B:180:GLU:HA	1:B:183:LEU:HD12	1.89	0.54
1:D:247:ARG:NE	1:D:256:GLU:OE2	2.38	0.54
1:C:119:LEU:HD23	1:C:127:LYS:HA	1.88	0.54
1:A:208:ALA:HA	1:A:237:LEU:HD22	1.90	0.54
1:B:309:THR:N	3:B:507:HOH:O	2.40	0.54
1:D:208:ALA:HA	1:D:237:LEU:HG	1.89	0.54
1:A:106:LEU:HD21	1:A:128:GLN:HG3	1.90	0.54
1:A:264:HIS:NE2	1:A:266:GLU:OE1	2.38	0.54
1:A:193:VAL:HG21	1:A:248:ALA:HB1	1.90	0.53
1:B:235:ASP:OD1	1:B:269:TYR:OH	2.25	0.53
1:A:388:TYR:HB3	2:E:8:PHE:HA	1.90	0.53
1:C:173:GLU:HG2	1:C:174:THR:H	1.74	0.53
1:C:199:LYS:HG3	1:C:244:LYS:HG2	1.91	0.53
1:D:240:GLN:CD	1:D:268:ARG:HH21	2.17	0.53
1:B:161:ASP:OD1	1:B:162:LEU:N	2.39	0.52
1:A:388:TYR:O	1:A:398:LEU:HA	2.09	0.52
2:F:4:GLN:O	2:F:9:TYR:OH	2.28	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:VAL:HG23	1:A:177:PHE:H	1.74	0.52
1:D:194:VAL:HG13	1:D:251:PRO:HA	1.91	0.52
1:B:120:ASP:HB3	1:B:126:LYS:HG2	1.91	0.52
1:C:193:VAL:HG21	1:C:248:ALA:HB1	1.91	0.52
1:C:169:ARG:NH1	1:C:171:SER:HB3	2.25	0.51
1:C:180:GLU:OE2	1:C:244:LYS:NZ	2.43	0.51
1:A:112:LEU:HD23	1:A:138:LEU:HB2	1.91	0.51
1:C:235:ASP:N	3:C:515:HOH:O	2.43	0.51
1:D:240:GLN:HE22	1:D:268:ARG:HE	1.58	0.51
1:D:199:LYS:HG3	1:D:244:LYS:HG2	1.92	0.51
2:G:4:GLN:HB3	2:G:5:PRO:HD2	1.91	0.51
1:B:235:ASP:OD2	1:B:235:ASP:N	2.43	0.51
1:C:207:SER:HG	1:C:212:LYS:H	1.58	0.51
1:D:112:LEU:HA	1:D:138:LEU:HD12	1.92	0.51
1:A:200:SER:O	1:A:242:THR:HA	2.11	0.51
1:A:206:LEU:HB2	1:A:237:LEU:HD21	1.93	0.51
1:C:133:VAL:HG23	1:C:135:SER:H	1.75	0.51
1:C:178:THR:N	1:C:181:SER:OG	2.43	0.50
1:C:313:VAL:HG12	1:C:354:ILE:HG22	1.92	0.50
1:C:247:ARG:NH1	1:C:256:GLU:OE1	2.45	0.50
1:A:107:VAL:HG12	1:A:109:ILE:HD11	1.92	0.50
1:A:350:LYS:HZ3	1:A:352:ILE:HD11	1.77	0.50
1:B:344:TRP:HA	1:B:354:ILE:H	1.76	0.50
1:A:350:LYS:NZ	1:A:352:ILE:HD11	2.28	0.49
1:B:325:LYS:NZ	3:B:510:HOH:O	2.45	0.49
1:B:236:ILE:HG13	1:C:252:ARG:HG3	1.93	0.49
1:A:344:TRP:N	3:A:507:HOH:O	2.46	0.49
1:B:312:LYS:HE3	1:B:357:PHE:CE2	2.48	0.49
1:D:116:ILE:O	1:D:131:LEU:N	2.45	0.49
1:B:198:GLY:HA3	1:C:200:SER:HA	1.95	0.49
1:D:182:LEU:HD23	1:D:195:LEU:HD13	1.93	0.48
1:B:106:LEU:HD13	1:B:118:ALA:HB1	1.95	0.48
1:A:350:LYS:NZ	3:A:514:HOH:O	2.46	0.48
1:B:112:LEU:HD23	1:B:392:TYR:HB2	1.95	0.48
1:C:300:GLU:N	1:C:300:GLU:OE2	2.47	0.48
1:D:375:ILE:HG13	1:D:408:PHE:HE1	1.79	0.48
2:G:4:GLN:HB3	2:G:5:PRO:CD	2.44	0.47
1:B:106:LEU:HA	1:B:121:PRO:HD3	1.97	0.47
1:A:395:GLN:NE2	1:A:396:LEU:O	2.48	0.47
1:C:314:SER:HB2	1:C:321:MET:HE2	1.96	0.47
1:D:318:TRP:CG	1:D:339:PRO:HA	2.49	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:7:ARG:HG3	2:F:8:PHE:H	1.79	0.47
1:B:315:VAL:HG12	1:B:357:PHE:O	2.14	0.47
1:A:180:GLU:HA	1:A:183:LEU:HD12	1.96	0.47
1:C:387:VAL:O	2:F:8:PHE:HA	2.15	0.47
1:D:105:SER:N	3:D:509:HOH:O	2.47	0.47
1:A:315:VAL:HG12	1:A:357:PHE:O	2.14	0.47
1:B:182:LEU:HB3	1:B:195:LEU:HD13	1.96	0.47
1:A:109:ILE:HD13	1:A:336:PHE:CZ	2.50	0.47
1:D:243:GLN:NE2	3:D:511:HOH:O	2.47	0.47
1:D:153:MET:HE1	2:G:3:PRO:HD3	1.96	0.47
1:D:224:GLN:OE1	1:D:224:GLN:N	2.48	0.46
1:D:375:ILE:HA	1:D:408:PHE:CE1	2.50	0.46
1:B:268:ARG:HH11	1:B:351:VAL:HB	1.80	0.46
1:A:237:LEU:H	1:A:237:LEU:HD23	1.80	0.46
1:D:107:VAL:HG23	1:D:121:PRO:HD3	1.96	0.46
1:D:236:ILE:HB	1:D:270:ILE:HB	1.97	0.46
1:C:165:TRP:HE3	1:C:172:MET:HB2	1.79	0.46
1:D:307:MET:O	1:D:309:THR:HG23	2.16	0.46
1:C:165:TRP:HE1	2:F:3:PRO:HD3	1.81	0.46
1:A:238:LEU:HD12	1:A:268:ARG:HB2	1.97	0.46
1:C:318:TRP:CE2	1:C:339:PRO:HG3	2.51	0.46
1:C:387:VAL:HG13	1:C:398:LEU:HD22	1.98	0.46
1:D:206:LEU:HD11	1:D:239:LEU:HD12	1.98	0.46
1:D:375:ILE:HG13	1:D:408:PHE:CE1	2.51	0.45
1:C:226:ASP:OD2	1:C:226:ASP:N	2.47	0.45
1:D:113:ASP:N	1:D:338:THR:HG21	2.30	0.45
1:D:240:GLN:NE2	1:D:268:ARG:HH21	2.14	0.45
1:D:400:SER:O	1:D:400:SER:OG	2.32	0.45
1:B:251:PRO:O	1:C:215:TYR:OH	2.21	0.45
1:A:247:ARG:NH1	1:A:256:GLU:OE2	2.49	0.45
1:A:318:TRP:CE2	1:A:339:PRO:HG3	2.51	0.45
1:A:389:LEU:HD11	1:A:396:LEU:HD23	1.99	0.45
1:A:358:ASP:HB3	2:E:8:PHE:CE2	2.52	0.45
1:D:141:SER:O	1:D:141:SER:OG	2.35	0.45
1:B:346:VAL:HG12	1:B:351:VAL:HG22	1.99	0.45
1:D:313:VAL:HG22	1:D:320:VAL:HG22	1.98	0.45
1:A:267:LEU:O	3:A:501:HOH:O	2.21	0.45
1:B:241:ARG:HG3	1:B:265:PHE:CZ	2.52	0.44
1:D:123:ASN:O	3:D:501:HOH:O	2.20	0.44
1:A:358:ASP:HB3	2:E:8:PHE:HE2	1.82	0.44
1:D:357:PHE:HB3	1:D:404:ILE:CG2	2.43	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:ARG:HA	1:B:223:ARG:HD2	1.90	0.44
1:D:318:TRP:CD1	1:D:318:TRP:N	2.84	0.44
1:A:238:LEU:HD13	1:D:194:VAL:HG11	1.99	0.44
1:B:108:ILE:HG21	1:B:267:LEU:HD13	1.99	0.44
1:D:389:LEU:HD11	1:D:396:LEU:HD22	1.99	0.44
1:A:109:ILE:HD13	1:A:336:PHE:HZ	1.83	0.43
1:D:214:ARG:HD3	1:D:214:ARG:HA	1.73	0.43
1:B:313:VAL:HG12	1:B:356:LEU:HA	2.01	0.43
1:B:313:VAL:HG11	1:B:356:LEU:HD22	1.99	0.43
1:C:304:ALA:HB1	1:C:345:LEU:HD23	2.00	0.43
1:D:324:ASN:ND2	1:D:326:LYS:O	2.47	0.43
1:A:199:LYS:HD3	1:A:201:LEU:HD11	1.99	0.43
1:A:251:PRO:HB2	1:D:270:ILE:HG21	2.01	0.43
1:C:269:TYR:HB2	1:C:351:VAL:HG13	2.01	0.43
1:C:313:VAL:HB	1:C:320:VAL:HG22	2.00	0.43
1:D:403:ARG:HD3	1:D:404:ILE:N	2.34	0.43
1:A:388:TYR:CD2	1:A:388:TYR:C	2.97	0.42
1:B:142:SER:O	1:B:258:TRP:HA	2.19	0.42
1:B:355:SER:HB2	1:B:357:PHE:CZ	2.54	0.42
1:C:387:VAL:HG22	1:C:398:LEU:HB3	2.00	0.42
1:A:331:GLU:OE1	3:A:503:HOH:O	2.21	0.42
1:C:389:LEU:HD12	1:C:397:TYR:O	2.19	0.42
1:D:147:GLU:N	3:D:502:HOH:O	2.52	0.42
1:D:143:LEU:HD23	1:D:144:SER:N	2.34	0.42
1:D:177:PHE:CD1	1:D:182:LEU:HD11	2.53	0.42
1:D:258:TRP:HB2	3:D:503:HOH:O	2.18	0.42
1:A:112:LEU:HD11	1:A:397:TYR:CZ	2.54	0.42
1:D:203:THR:HG22	1:D:217:CYS:HB2	2.02	0.42
1:D:388:TYR:HB2	2:G:8:PHE:CE1	2.54	0.42
1:C:112:LEU:HD11	1:C:397:TYR:CZ	2.54	0.42
1:C:161:ASP:N	1:C:161:ASP:OD1	2.52	0.42
1:C:207:SER:OG	1:C:212:LYS:N	2.47	0.42
1:D:408:PHE:N	3:D:507:HOH:O	2.50	0.42
1:A:238:LEU:HD11	1:A:270:ILE:HD11	2.01	0.42
1:A:311:ILE:HA	1:A:322:ALA:HA	2.02	0.42
1:B:133:VAL:HG21	1:B:265:PHE:HZ	1.84	0.42
1:B:106:LEU:O	1:B:346:VAL:HG22	2.19	0.42
1:B:203:THR:OG1	1:B:240:GLN:OE1	2.34	0.41
1:B:324:ASN:OD1	1:B:326:LYS:N	2.48	0.41
1:D:112:LEU:C	1:D:114:GLY:H	2.28	0.41
1:C:319:LYS:HZ2	1:C:333:GLU:CD	2.28	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:240:GLN:N	1:D:240:GLN:OE1	2.52	0.41
1:D:108:ILE:N	1:D:344:TRP:O	2.53	0.41
1:B:164:GLN:O	1:B:173:GLU:HG3	2.21	0.41
2:F:4:GLN:HA	2:F:5:PRO:HD3	1.82	0.41
1:B:243:GLN:NE2	1:B:261:SER:OG	2.54	0.41
1:B:196:VAL:O	1:B:247:ARG:N	2.52	0.41
1:C:238:LEU:HD23	1:C:238:LEU:HA	1.79	0.41
1:D:104:ARG:N	3:D:509:HOH:O	2.53	0.41
1:B:208:ALA:HA	1:B:237:LEU:HD12	2.02	0.41
1:B:252:ARG:NH2	1:C:235:ASP:OD2	2.54	0.41
1:A:238:LEU:HD12	1:A:268:ARG:HE	1.86	0.41
1:C:152:LYS:NZ	3:C:517:HOH:O	2.45	0.41
1:C:154:ILE:HD13	1:C:164:GLN:HB2	2.02	0.41
1:A:164:GLN:HG2	1:A:173:GLU:O	2.21	0.41
1:A:179:VAL:HG11	1:A:244:LYS:HB3	2.02	0.41
1:A:324:ASN:HB2	1:A:331:GLU:HG2	2.03	0.41
1:C:358:ASP:O	1:C:404:ILE:HG22	2.21	0.41
1:A:113:ASP:H	1:A:338:THR:HG21	1.87	0.40
1:A:346:VAL:HA	1:A:350:LYS:O	2.21	0.40
2:E:7:ARG:NH2	3:E:101:HOH:O	2.40	0.40
1:D:360:THR:HG22	1:D:404:ILE:O	2.22	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	179/326 (55%)	168 (94%)	11 (6%)	0	100	100
1	B	172/326 (53%)	163 (95%)	9 (5%)	0	100	100
1	C	232/326 (71%)	217 (94%)	15 (6%)	0	100	100
1	D	235/326 (72%)	210 (89%)	23 (10%)	2 (1%)	14	41

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	7/12 (58%)	4 (57%)	3 (43%)	0	100	100
2	F	7/12 (58%)	5 (71%)	1 (14%)	1 (14%)	0	0
2	G	7/12 (58%)	4 (57%)	2 (29%)	1 (14%)	0	0
All	All	839/1340 (63%)	771 (92%)	64 (8%)	4 (0%)	24	55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	G	5	PRO
2	F	5	PRO
1	D	145	LYS
1	D	376	VAL

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	166/285 (58%)	162 (98%)	4 (2%)	43	77
1	B	164/285 (58%)	160 (98%)	4 (2%)	43	77
1	C	212/285 (74%)	207 (98%)	5 (2%)	43	77
1	D	215/285 (75%)	207 (96%)	8 (4%)	30	65
2	E	8/10 (80%)	8 (100%)	0	100	100
2	F	8/10 (80%)	8 (100%)	0	100	100
2	G	8/10 (80%)	8 (100%)	0	100	100
All	All	781/1170 (67%)	760 (97%)	21 (3%)	39	74

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	155	ILE
1	A	330	LEU
1	A	357	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	388	TYR
1	B	119	LEU
1	B	194	VAL
1	B	346	VAL
1	B	356	LEU
1	C	184	GLU
1	C	193	VAL
1	C	313	VAL
1	C	324	ASN
1	C	407	LYS
1	D	116	ILE
1	D	245	THR
1	D	246	VAL
1	D	262	VAL
1	D	301	GLU
1	D	377	GLU
1	D	380	ARG
1	D	408	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	GLN
1	A	335	GLN
1	B	255	ASN
1	C	240	GLN
1	C	329	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	193/326 (59%)	0.71	12 (6%) 26 20	43, 58, 70, 83	0
1	B	190/326 (58%)	0.68	9 (4%) 36 29	43, 61, 79, 100	0
1	C	244/326 (74%)	0.56	7 (2%) 53 43	34, 50, 65, 71	0
1	D	247/326 (75%)	0.48	6 (2%) 59 49	37, 49, 64, 79	0
2	E	9/12 (75%)	1.13	0 100 100	71, 76, 82, 84	0
2	F	9/12 (75%)	1.17	2 (22%) 2 1	63, 68, 77, 87	0
2	G	9/12 (75%)	1.41	2 (22%) 2 1	51, 59, 70, 73	0
All	All	901/1340 (67%)	0.62	38 (4%) 40 32	34, 54, 73, 100	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	324	ASN	4.2
1	C	159	ASP	4.0
1	B	135	SER	3.9
2	F	4	GLN	3.8
1	B	205	GLY	3.6
1	C	381	GLY	3.6
1	A	390	GLY	3.2
1	A	249	VAL	2.8
1	C	401	SER	2.7
1	A	311	ILE	2.7
1	A	322	ALA	2.6
1	D	108	ILE	2.6
2	G	9	TYR	2.6
1	A	323	PHE	2.6
2	G	7	ARG	2.6
1	B	389	LEU	2.5
1	B	130	ASP	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	131	LEU	2.5
1	D	198	GLY	2.5
1	A	134	GLY	2.4
1	D	304	ALA	2.4
1	B	269	TYR	2.4
1	D	360	THR	2.4
1	A	211	GLY	2.3
1	C	198	GLY	2.3
1	A	329	HIS	2.3
1	C	271	PRO	2.2
1	C	263	GLY	2.2
1	D	135	SER	2.2
1	C	382	ALA	2.2
1	B	106	LEU	2.1
1	D	110	SER	2.1
2	F	8	PHE	2.1
1	A	133	VAL	2.1
1	A	141	SER	2.0
1	B	270	ILE	2.0
1	A	242	THR	2.0
1	B	204	TYR	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.