



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 08:42 PM UTC

PDB ID : 1CP3 / pdb_00001cp3
Title : CRYSTAL STRUCTURE OF THE COMPLEX OF APOPAIN WITH THE
TETRAPEPTIDE INHIBITOR ACE-DVAD-FMC
Authors : Mittl, P.R.E.; Dimarco, S.; Gruetter, M.G.
Deposited on : 1996-12-12
Resolution : 2.30 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

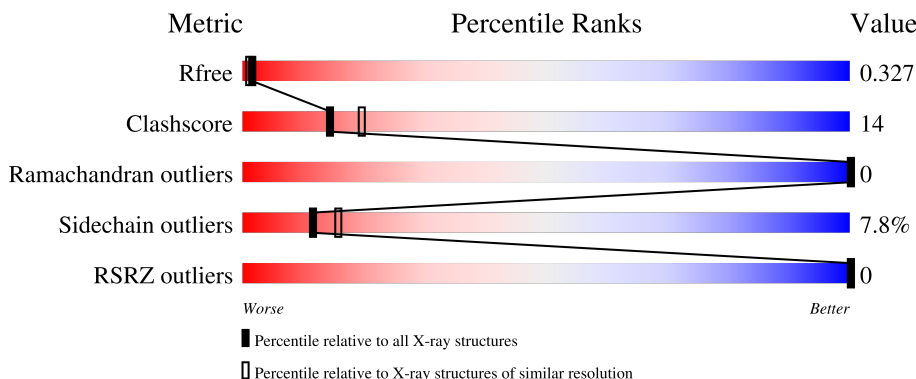
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

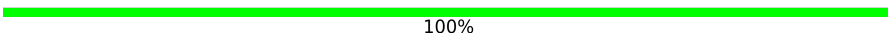
The reported resolution of this entry is 2.30 Å.

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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	277	
1	B	277	
2	C	6	
2	D	6	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4174 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 APOPAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	0	0
			1882	1194	325	349	14			
1	B	232	Total	C	N	O	S	0	0	0
			1882	1194	325	349	14			

- Molecule 2 is a protein called ACETYL-ASP-VAL-ALA-ASP-FLUOROMETHYLKETON E.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	0	1
			32	19	4	9			
2	D	6	Total	C	N	O	0	0	1
			32	19	4	9			

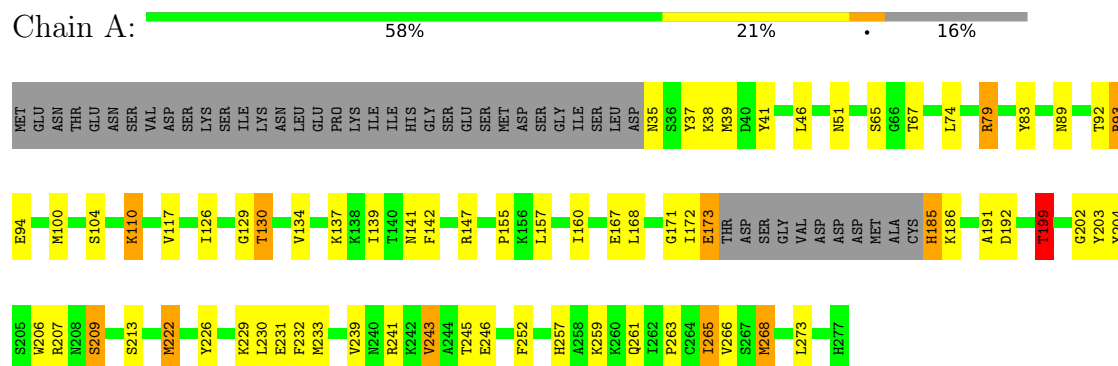
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	200	Total	O	0	0
			200	200		
3	C	6	Total	O	0	0
			6	6		
3	B	136	Total	O	0	0
			136	136		
3	D	4	Total	O	0	0
			4	4		

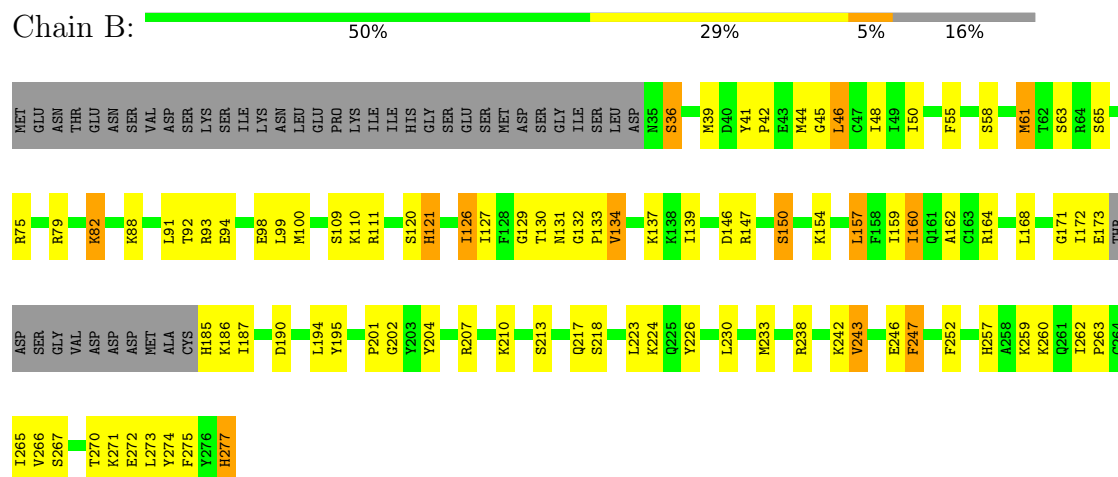
3 Residue-property plots

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: APOPAIN



• Molecule 1: APOPAIN



• Molecule 2: ACETYL-ASP-VAL-ALA-ASP-FLUOROMETHYLKETONE



There are no outlier residues recorded for this chain.

• Molecule 2: ACETYL-ASP-VAL-ALA-ASP-FLUOROMETHYLKETONE



ACE889
D990
V991
A992
D993
CF0994

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.90Å 69.10Å 93.80Å 90.00° 101.20° 90.00°	Depositor
Resolution (Å)	8.00 – 2.30 8.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	83.7 (8.00-2.30) 79.4 (8.00-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 2.31Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.188 , 0.284 0.312 , 0.327	Depositor DCC
R_{free} test set	2383 reflections (9.90%)	wwPDB-VP
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4174	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CF0, ACE

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	0.82	1/1922 (0.1%)	1.12	15/2581 (0.6%)
1	B	0.72	0/1922	1.12	16/2581 (0.6%)
2	C	0.72	0/28	0.69	0/38
2	D	0.65	0/28	0.95	0/38
All	All	0.77	1/3900 (0.0%)	1.11	31/5238 (0.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	268	MET	SD-CE	-5.21	1.66	1.79

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	129	GLY	N-CA-C	-10.01	97.28	112.60
1	A	129	GLY	N-CA-C	-8.36	99.12	112.66
1	A	226	TYR	N-CA-C	7.97	124.08	113.72
1	B	160	ILE	N-CA-C	6.63	117.58	107.77
1	B	126	ILE	N-CA-C	6.53	118.18	108.46
1	B	133	PRO	N-CA-C	6.38	120.44	110.80
1	B	267	SER	N-CA-C	6.31	119.53	109.24
1	A	160	ILE	N-CA-C	6.21	116.81	107.75
1	B	226	TYR	N-CA-C	6.04	122.68	113.61
1	A	268	MET	N-CA-C	-5.99	104.62	112.41
1	A	265	ILE	CB-CA-C	-5.88	103.62	110.91
1	A	199	THR	CB-CA-C	-5.83	97.85	111.09
1	B	132	GLY	CA-C-N	5.71	125.90	119.90
1	B	132	GLY	C-N-CA	5.71	125.90	119.90
1	A	246	GLU	N-CA-C	5.62	120.19	113.23
1	A	230	LEU	N-CA-C	5.59	119.08	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	203	TYR	N-CA-C	5.57	119.05	110.42
1	A	157	LEU	N-CA-C	5.51	117.71	108.73
1	A	231	GLU	N-CA-C	-5.50	102.33	110.48
1	B	246	GLU	N-CA-C	5.39	119.92	113.23
1	B	109	SER	N-CA-C	5.36	117.54	111.11
1	B	79	ARG	N-CA-C	-5.30	105.55	112.23
1	B	134	VAL	CB-CA-C	-5.29	102.78	110.81
1	A	206	TRP	N-CA-C	5.23	118.59	110.17
1	B	252	PHE	N-CA-C	-5.21	106.76	113.23
1	B	154	LYS	CA-C-N	5.19	125.35	119.90
1	B	154	LYS	C-N-CA	5.19	125.35	119.90
1	B	121	HIS	N-CA-C	-5.13	102.93	110.52
1	A	186	LYS	N-CA-C	5.06	117.64	110.50
1	A	79	ARG	N-CA-C	-5.05	105.48	111.69
1	A	232	PHE	N-CA-C	5.03	117.15	111.11

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	1882	0	1854	48	3
1	B	1882	0	1854	68	0
2	C	32	0	25	0	0
2	D	32	0	25	2	0
3	A	200	0	0	7	3
3	B	136	0	0	3	0
3	C	6	0	0	0	0
3	D	4	0	0	0	0
All	All	4174	0	3758	107	5

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 14.

All (107) 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:A:233:MET:HE1	1:A:265:ILE:HG21	1.62	0.80
1:A:100:MET:HG3	1:A:139:ILE:HG23	1.63	0.80
1:B:150:SER:HB2	3:B:350:HOH:O	1.87	0.73
1:B:46:LEU:HD23	1:B:111:ARG:NH1	2.05	0.71
1:A:243:VAL:HG22	1:A:263:PRO:HD3	1.73	0.71
1:A:110:LYS:H	1:A:110:LYS:HZ2	1.37	0.69
1:A:222:MET:HE3	1:A:222:MET:HA	1.73	0.69
1:A:199:THR:HG21	1:A:204:TYR:O	1.93	0.67
1:A:233:MET:HE2	1:B:233:MET:HE2	1.76	0.67
1:B:50:ILE:HG23	1:B:91:LEU:HD23	1.77	0.66
1:B:233:MET:HE1	1:B:265:ILE:HG21	1.78	0.66
1:A:199:THR:HG22	3:A:345:HOH:O	1.95	0.65
1:B:61:MET:HA	1:B:61:MET:HE3	1.77	0.65
1:B:46:LEU:HD23	1:B:111:ARG:CZ	2.28	0.64
1:B:274:TYR:HB2	1:B:277:HIS:O	1.98	0.64
1:A:243:VAL:CG2	1:A:263:PRO:HD3	2.28	0.63
1:B:92:THR:HG22	1:B:130:THR:OG1	1.98	0.63
1:A:222:MET:HA	1:A:222:MET:CE	2.30	0.62
1:A:172:ILE:O	1:B:186:LYS:HA	2.00	0.61
1:B:93:ARG:HG3	1:B:134:VAL:HG22	1.82	0.61
1:B:39:MET:HE2	1:B:273:LEU:HD11	1.84	0.59
1:B:100:MET:HG3	1:B:139:ILE:HG23	1.86	0.58
1:B:257:HIS:O	1:B:259:LYS:HE3	2.03	0.58
1:B:164:ARG:O	1:B:202:GLY:HA2	2.03	0.57
1:A:233:MET:HE2	1:B:233:MET:CE	2.33	0.57
1:B:82:LYS:N	1:B:82:LYS:HD3	2.20	0.56
1:B:185:HIS:HB3	3:B:326:HOH:O	2.06	0.55
1:B:39:MET:HE3	1:B:44:MET:HE3	1.88	0.55
1:A:130:THR:HG21	3:A:348:HOH:O	2.05	0.55
1:B:157:LEU:HD12	1:B:194:LEU:HB3	1.87	0.55
1:B:172:ILE:HD12	1:B:173:GLU:H	1.71	0.55
1:B:213:SER:O	1:B:217:GLN:HG3	2.06	0.55
1:A:239:VAL:O	1:A:243:VAL:HG13	2.08	0.54
1:B:195:TYR:HB2	1:B:266:VAL:HB	1.91	0.53
1:A:141:ASN:HB3	1:A:147:ARG:NH1	2.23	0.52
1:A:245:THR:HG22	1:B:185:HIS:HE1	1.75	0.52
1:B:233:MET:HE1	1:B:265:ILE:CG2	2.39	0.52
1:A:93:ARG:HB2	1:A:134:VAL:HG22	1.91	0.52
1:A:207:ARG:HA	1:A:213:SER:HA	1.92	0.52
1:A:199:THR:HG23	1:A:261:GLN:HG3	1.91	0.51
1:B:36:SER:HA	1:B:271:LYS:HB3	1.92	0.51
1:B:243:VAL:HG22	1:B:263:PRO:HD3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:LYS:HB3	1:B:91:LEU:HD21	1.92	0.51
1:B:207:ARG:NH2	2:D:991:VAL:HG21	2.26	0.51
1:B:61:MET:HG3	1:B:121:HIS:CD2	2.46	0.50
1:A:35:ASN:N	1:A:35:ASN:OD1	2.45	0.50
1:A:185:HIS:HB2	3:A:457:HOH:O	2.12	0.50
1:A:74:LEU:HD13	1:A:117:VAL:HG11	1.93	0.49
1:B:218:SER:HB3	1:B:242:LYS:HD3	1.94	0.49
1:B:247:PHE:CD1	1:B:247:PHE:N	2.80	0.49
1:B:172:ILE:O	1:B:173:GLU:HB3	2.12	0.49
1:A:233:MET:CE	1:B:233:MET:HE2	2.41	0.49
1:B:46:LEU:CD1	1:B:48:ILE:HD11	2.42	0.49
1:B:92:THR:HG22	1:B:130:THR:HG1	1.76	0.49
1:B:147:ARG:HH11	1:B:147:ARG:HG3	1.78	0.48
1:A:92:THR:HG22	1:A:130:THR:HG22	1.95	0.48
1:A:167:GLU:HB2	1:A:202:GLY:O	2.14	0.47
1:B:224:LYS:HB3	1:B:224:LYS:HE3	1.70	0.47
1:A:110:LYS:NZ	3:A:321:HOH:O	2.45	0.47
1:B:164:ARG:NH1	1:B:201:PRO:HG3	2.29	0.47
1:A:39:MET:HE2	1:A:273:LEU:HD11	1.97	0.47
1:B:218:SER:CB	1:B:242:LYS:HD3	2.45	0.46
1:B:207:ARG:HA	1:B:213:SER:HA	1.98	0.46
1:B:262:ILE:HD12	1:B:263:PRO:HD2	1.97	0.46
1:A:233:MET:CE	1:A:265:ILE:HG21	2.40	0.46
1:B:36:SER:HB3	1:B:272:GLU:HB2	1.98	0.46
1:A:168:LEU:HD22	1:A:259:LYS:HG3	1.98	0.46
1:A:172:ILE:HG13	1:A:173:GLU:N	2.30	0.46
1:B:217:GLN:HG2	3:B:333:HOH:O	2.15	0.46
1:B:41:TYR:HB3	1:B:42:PRO:HD2	1.97	0.45
1:B:92:THR:HB	1:B:131:ASN:CG	2.42	0.45
1:A:222:MET:HE3	1:A:222:MET:CA	2.44	0.45
1:B:120:SER:OG	1:B:121:HIS:N	2.50	0.45
1:B:162:ALA:HA	2:D:993:ASP:HB3	1.98	0.45
1:A:137:LYS:HE3	3:A:399:HOH:O	2.18	0.44
1:A:38:LYS:HD3	1:A:41:TYR:CE1	2.52	0.44
1:B:168:LEU:CD2	1:B:204:TYR:CD2	3.01	0.44
1:B:110:LYS:HD2	1:B:110:LYS:H	1.83	0.43
1:B:44:MET:HE1	1:B:275:PHE:O	2.17	0.43
1:A:172:ILE:HD13	1:B:146:ASP:HA	1.99	0.43
1:B:210:LYS:HE3	1:B:210:LYS:HB2	1.88	0.43
1:B:55:PHE:HB2	1:B:61:MET:O	2.18	0.43
1:B:50:ILE:HD11	1:B:99:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ASN:N	3:A:391:HOH:O	2.51	0.42
1:A:92:THR:HA	1:A:130:THR:HG22	2.01	0.42
1:B:243:VAL:HA	1:B:247:PHE:CD1	2.55	0.42
1:A:51:ASN:ND2	1:A:67:THR:OG1	2.53	0.42
1:A:37:TYR:HA	3:A:357:HOH:O	2.19	0.41
1:A:51:ASN:HB3	1:A:89:ASN:ND2	2.35	0.41
1:B:39:MET:HB2	1:B:277:HIS:OXT	2.20	0.41
1:B:168:LEU:CD2	1:B:204:TYR:HD2	2.32	0.41
1:B:172:ILE:HG13	1:B:173:GLU:N	2.35	0.41
1:A:191:ALA:O	1:A:192:ASP:HB2	2.20	0.41
1:A:65:SER:O	1:A:209:SER:HA	2.20	0.41
1:A:171:GLY:HA2	1:B:187:ILE:O	2.21	0.41
1:A:241:ARG:HB2	1:B:270:THR:O	2.20	0.41
1:B:45:GLY:HA2	1:B:111:ARG:HB3	2.02	0.41
1:A:37:TYR:CD1	1:A:155:PRO:HD3	2.55	0.41
1:B:137:LYS:HA	1:B:137:LYS:HD3	1.84	0.41
1:A:266:VAL:HA	1:B:265:ILE:O	2.21	0.41
1:B:164:ARG:HH12	1:B:201:PRO:HG3	1.85	0.41
1:B:171:GLY:HA3	1:B:260:LYS:HG3	2.03	0.41
1:A:141:ASN:HB3	1:A:147:ARG:HH11	1.87	0.40
1:A:233:MET:HE1	1:A:265:ILE:CG2	2.40	0.40
1:A:104:SER:OG	1:A:142:PHE:HB3	2.21	0.40
1:B:230:LEU:HD11	1:B:238:ARG:NH1	2.35	0.40
1:B:159:ILE:C	1:B:160:ILE:HG13	2.46	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:HIS:NE2	3:A:333:HOH:O[2_656]	1.38	0.82
1:A:83:TYR:O	1:A:252:PHE:CE1[2_646]	1.80	0.40
3:A:403:HOH:O	3:A:437:HOH:O[2_646]	1.82	0.38
3:A:410:HOH:O	3:A:427:HOH:O[2_646]	1.84	0.36
1:A:79:ARG:NH1	1:A:252:PHE:O[2_646]	1.96	0.24

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	228/277 (82%)	220 (96%)	8 (4%)	0	100	100
1	B	228/277 (82%)	216 (95%)	12 (5%)	0	100	100
2	C	3/6 (50%)	3 (100%)	0	0	100	100
2	D	3/6 (50%)	3 (100%)	0	0	100	100
All	All	462/566 (82%)	442 (96%)	20 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	208/249 (84%)	194 (93%)	14 (7%)	15	21
1	B	208/249 (84%)	189 (91%)	19 (9%)	9	11
2	C	3/3 (100%)	3 (100%)	0	100	100
2	D	3/3 (100%)	3 (100%)	0	100	100
All	All	422/504 (84%)	389 (92%)	33 (8%)	11	16

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU
1	A	93	ARG

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Mol	Chain	Res	Type
1	A	94	GLU
1	A	110	LYS
1	A	126	ILE
1	A	130	THR
1	A	173	GLU
1	A	185	HIS
1	A	199	THR
1	A	209	SER
1	A	222	MET
1	A	229	LYS
1	A	243	VAL
1	A	268	MET
1	B	36	SER
1	B	46	LEU
1	B	58	SER
1	B	61	MET
1	B	63	SER
1	B	65	SER
1	B	75	ARG
1	B	82	LYS
1	B	94	GLU
1	B	98	GLU
1	B	126	ILE
1	B	127	ILE
1	B	150	SER
1	B	157	LEU
1	B	190	ASP
1	B	223	LEU
1	B	243	VAL
1	B	247	PHE
1	B	277	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	51	ASN
1	A	89	ASN
1	A	217	GLN
1	B	51	ASN
1	B	56	HIS
1	B	89	ASN

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Mol	Chain	Res	Type
1	B	121	HIS
1	B	141	ASN
1	B	161	GLN
1	B	185	HIS

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](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	232/277 (83%)	-1.41	0 100 100	8, 22, 42, 73	0
1	B	232/277 (83%)	-1.37	0 100 100	14, 34, 60, 96	0
2	C	4/6 (66%)	-1.50	0 100 100	19, 20, 27, 27	0
2	D	4/6 (66%)	-1.59	0 100 100	29, 31, 35, 36	0
All	All	472/566 (83%)	-1.39	0 100 100	8, 28, 54, 96	0

There are no RSRZ outliers to report.

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.