



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:40 AM GMT

PDB ID : 3BPF
Title : Crystal Structure of Falcipain-2 with Its inhibitor, E64
Authors : kerr, I.D.; Lee, J.H.; Brinen, L.S.
Deposited on : 2007-12-18
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

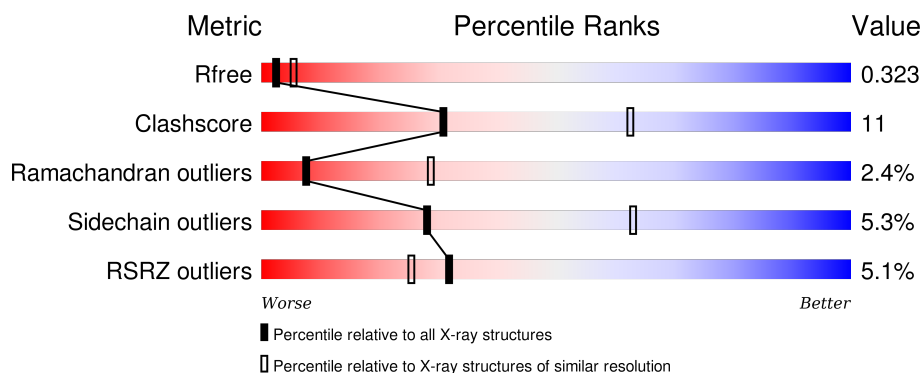
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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. 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	241	<div> <div>4%</div> <div>67% 25% 6% •</div> </div>
1	B	241	<div> <div>5%</div> <div>77% 21% •</div> </div>
1	C	241	<div> <div>6%</div> <div>77% 20% • •</div> </div>
1	D	241	<div> <div>5%</div> <div>67% 21% • 11%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7453 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.

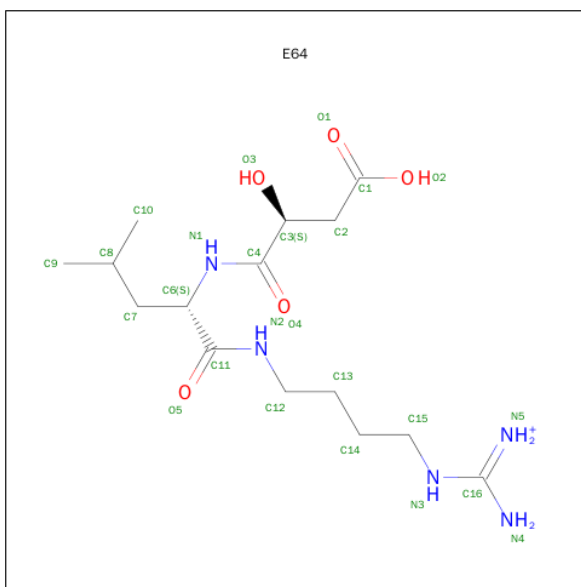
- Molecule 1 is a protein called Cysteine protease falcipain-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	236	Total	C	N	O	S	0	0	0
			1868	1183	308	363	14			
1	B	240	Total	C	N	O	S	0	0	0
			1898	1205	314	365	14			
1	C	234	Total	C	N	O	S	0	0	0
			1847	1174	303	356	14			
1	D	215	Total	C	N	O	S	0	0	0
			1692	1076	275	328	13			

There are 4 discrepancies between the modelled and reference sequences:

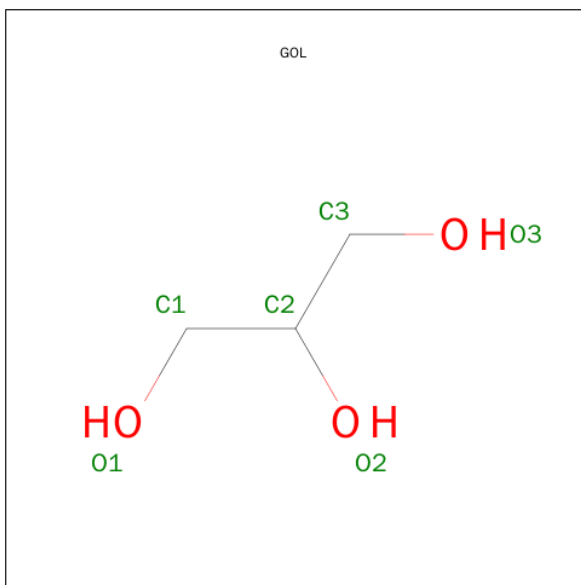
Chain	Residue	Modelled	Actual	Comment	Reference
A	171	GLN	GLU	SEE REMARK 999	UNP Q9NBD4
B	171	GLN	GLU	SEE REMARK 999	UNP Q9NBD4
C	171	GLN	GLU	SEE REMARK 999	UNP Q9NBD4
D	171	GLN	GLU	SEE REMARK 999	UNP Q9NBD4

- Molecule 2 is N-[N-[1-HYDROXYCARBOXYETHYL-CARBONYL]LEUCYLAMINO-BUTYL]-GUANIDINE (three-letter code: E64) (formula: C₁₅H₃₀N₅O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			25	15	5	5		
2	B	1	Total	C	N	O	0	0
			25	15	5	5		
2	C	1	Total	C	N	O	0	0
			25	15	5	5		
2	D	1	Total	C	N	O	0	0
			25	15	5	5		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			6	3	3		

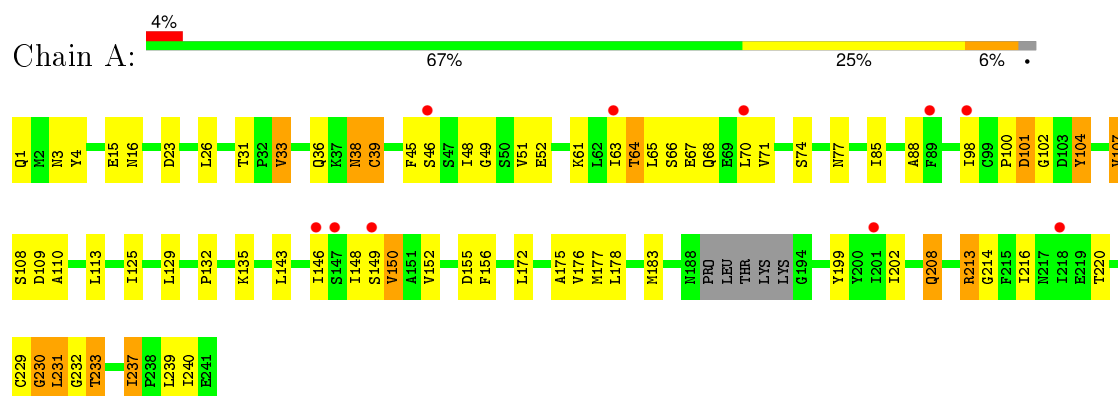
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	10	Total	O	0	0
			10	10		
4	B	13	Total	O	0	0
			13	13		
4	C	15	Total	O	0	0
			15	15		
4	D	4	Total	O	0	0
			4	4		

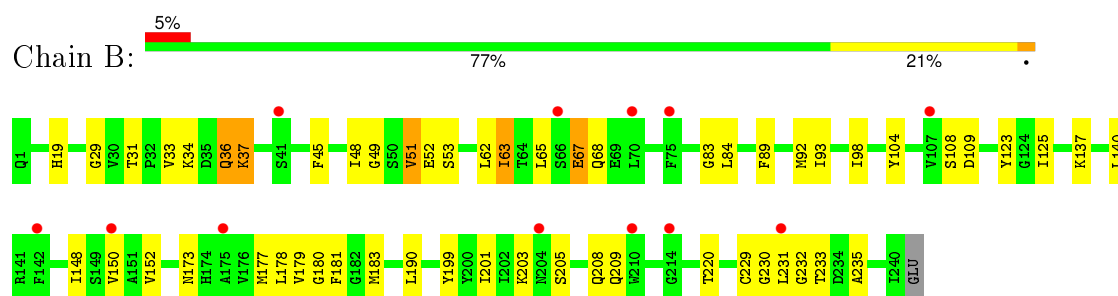
3 Residue-property plots [i](#)

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.

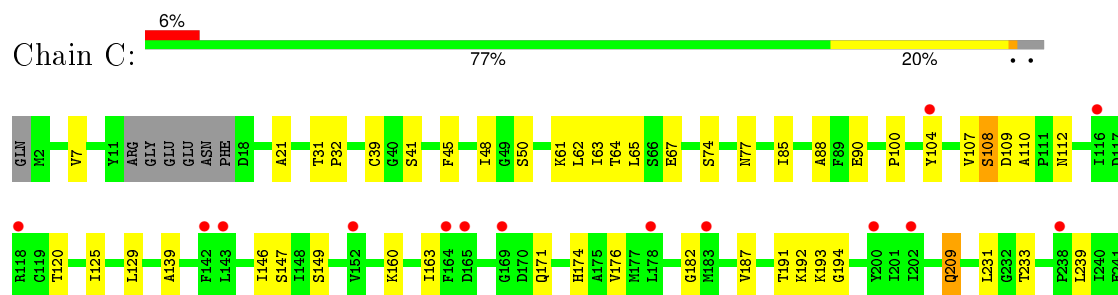
• Molecule 1: Cysteine protease falcipain-2



• Molecule 1: Cysteine protease falcipain-2

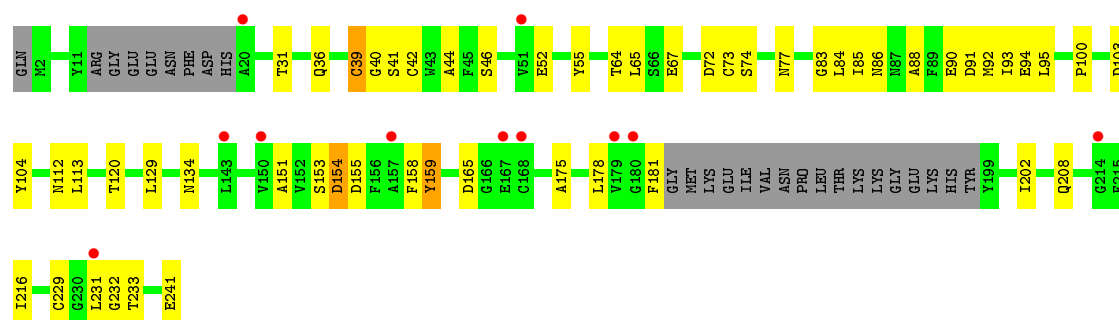


• Molecule 1: Cysteine protease falcipain-2



• Molecule 1: Cysteine protease falcipain-2





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	143.78Å 167.81Å 177.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.67 – 2.90 66.65 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.9 (66.67-2.90) 99.9 (66.65-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.275 , 0.325 0.275 , 0.323	Depositor DCC
R_{free} test set	2419 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	69.8	Xtriage
Anisotropy	0.189	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 107.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 52560 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	7453	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.86% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, E64

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	0.46	0/1909	0.58	0/2572
1	B	0.37	0/1941	0.51	0/2618
1	C	0.39	0/1888	0.50	0/2545
1	D	0.40	0/1728	0.51	0/2330
All	All	0.41	0/7466	0.53	0/10065

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	36	GLN	Peptide

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	1868	0	1774	56	0
1	B	1898	0	1820	32	0
1	C	1847	0	1773	29	0
1	D	1692	0	1613	22	0
2	A	25	0	28	1	0
2	B	25	0	28	0	0
2	C	25	0	28	1	0
2	D	25	0	28	1	0
3	C	6	0	8	0	0
4	A	10	0	0	0	0
4	B	13	0	0	0	0
4	C	15	0	0	0	0
4	D	4	0	0	0	0
All	All	7453	0	7100	137	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 11.

All (137) 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:67:GLU:O	1:A:71:VAL:HG23	1.63	0.97
1:C:129:LEU:HD23	1:C:239:LEU:HD11	1.41	0.97
1:A:129:LEU:HD22	1:A:239:LEU:HD11	1.61	0.82
1:C:174:HIS:CD2	1:C:176:VAL:HG13	2.21	0.76
1:D:83:GLY:O	1:D:84:LEU:HD23	1.87	0.75
1:A:129:LEU:CD2	1:A:237:ILE:HG23	2.21	0.70
1:C:7:VAL:HG21	1:C:129:LEU:HD22	1.75	0.68
1:A:129:LEU:HD21	1:A:237:ILE:HG23	1.75	0.68
1:C:108:SER:OG	1:C:109:ASP:N	2.21	0.68
1:B:63:ILE:HG21	1:B:123:TYR:CE1	2.29	0.68
1:C:21:ALA:HB2	1:C:182:GLY:HA2	1.77	0.67
1:B:48:ILE:HD12	1:B:49:GLY:N	2.11	0.66
1:A:23:ASP:HB3	1:A:26:LEU:HD12	1.77	0.65
1:B:199:TYR:HA	1:B:220:THR:HG22	1.78	0.65
1:A:178:LEU:HD23	1:A:202:ILE:HD13	1.79	0.64
1:A:113:LEU:HD21	1:C:31:THR:HA	1.79	0.64
1:D:72:ASP:OD2	1:D:112:ASN:ND2	2.31	0.63
1:B:33:VAL:HG13	1:B:208:GLN:HG2	1.80	0.63
1:A:178:LEU:HD23	1:A:202:ILE:CD1	2.28	0.62
1:A:143:LEU:HB3	1:A:237:ILE:HD13	1.80	0.62
1:C:107:VAL:HG12	1:C:110:ALA:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:LYS:NZ	1:C:63:ILE:HD11	2.14	0.62
1:A:107:VAL:O	1:A:107:VAL:HG23	2.01	0.61
1:A:48:ILE:HD11	1:A:98:ILE:HD13	1.83	0.61
1:C:21:ALA:CB	1:C:182:GLY:HA2	2.31	0.59
1:C:174:HIS:HD2	1:C:176:VAL:HG13	1.68	0.58
1:B:33:VAL:CG1	1:B:208:GLN:HG2	2.34	0.58
1:D:151:ALA:HB3	1:D:229:CYS:HB3	1.86	0.58
1:A:199:TYR:HA	1:A:220:THR:HG22	1.85	0.58
1:A:33:VAL:HG11	1:A:214:GLY:HA3	1.86	0.57
1:B:177:MET:HE3	1:B:203:LYS:HD3	1.85	0.57
1:B:148:ILE:HD11	1:B:178:LEU:HD13	1.85	0.57
1:A:31:THR:HG22	1:D:113:LEU:HD21	1.86	0.56
1:B:29:GLY:O	1:B:53:SER:OG	2.16	0.56
1:A:150:VAL:HG23	1:A:229:CYS:O	2.06	0.55
1:A:202:ILE:HB	1:A:216:ILE:HG23	1.88	0.54
1:D:39:CYS:O	1:D:41:SER:N	2.40	0.54
1:B:150:VAL:HG23	1:B:229:CYS:HB2	1.91	0.53
1:B:33:VAL:HG22	1:B:34:LYS:O	2.08	0.53
1:D:85:ILE:O	1:D:88:ALA:HB3	2.09	0.53
1:A:107:VAL:HG23	1:A:110:ALA:HB3	1.91	0.53
1:B:108:SER:OG	1:B:109:ASP:N	2.35	0.52
1:D:86:ASN:N	1:D:86:ASN:OD1	2.40	0.52
1:A:108:SER:OG	1:A:109:ASP:N	2.41	0.52
1:D:52:GLU:OE2	1:D:64:THR:HA	2.09	0.52
1:A:208:GLN:HB3	1:A:213:ARG:HG2	1.90	0.52
1:A:70:LEU:HD12	1:A:74:SER:HB2	1.92	0.52
1:A:149:SER:HB2	1:A:172:LEU:HD22	1.92	0.52
1:A:4:TYR:HA	1:A:129:LEU:HD12	1.92	0.51
1:A:101:ASP:OD1	1:A:102:GLY:N	2.44	0.51
1:B:36:GLN:HG2	1:B:205:SER:O	2.10	0.51
1:A:74:SER:OG	1:A:77:ASN:ND2	2.41	0.51
1:C:39:CYS:SG	1:C:41:SER:OG	2.69	0.51
1:D:44:ALA:HB1	1:D:67:GLU:HB2	1.93	0.51
1:A:199:TYR:CA	1:A:220:THR:HG22	2.40	0.51
1:B:19:HIS:O	1:B:137:LYS:NZ	2.38	0.50
1:A:46:SER:CB	1:A:175:ALA:HB1	2.42	0.50
1:A:33:VAL:HG22	1:A:208:GLN:HG2	1.94	0.50
1:D:88:ALA:O	1:D:92:MET:HG3	2.12	0.50
1:B:140:LEU:HD11	1:B:179:VAL:C	2.32	0.50
1:C:74:SER:OG	1:C:77:ASN:ND2	2.39	0.49
1:A:176:VAL:CG2	1:A:202:ILE:HD12	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:139:ALA:CB	1:C:146:ILE:HD11	2.42	0.49
1:C:107:VAL:CG1	1:C:110:ALA:HB3	2.42	0.49
1:A:156:PHE:CZ	1:A:216:ILE:HD13	2.47	0.49
1:D:93:ILE:HG22	1:D:94:GLU:N	2.26	0.49
1:C:50:SER:HB2	1:C:147:SER:OG	2.12	0.49
1:B:181:PHE:HA	1:B:201:ILE:HD12	1.94	0.49
1:A:150:VAL:HG23	1:A:229:CYS:HB2	1.94	0.48
1:B:48:ILE:HA	1:B:92:MET:HE1	1.95	0.48
1:C:45:PHE:CE2	1:C:67:GLU:HG3	2.49	0.48
1:C:191:THR:O	1:C:193:LYS:N	2.47	0.48
1:A:85:ILE:O	1:A:88:ALA:HB3	2.13	0.47
1:A:176:VAL:HG21	1:A:202:ILE:HD12	1.96	0.47
1:B:148:ILE:HG22	1:B:235:ALA:HA	1.95	0.47
1:A:129:LEU:HD23	1:A:237:ILE:HG23	1.96	0.47
1:A:172:LEU:HD11	1:A:233:THR:OG1	2.14	0.47
1:D:91:ASP:O	1:D:95:LEU:HD12	2.15	0.46
1:C:31:THR:HB	1:C:32:PRO:HD2	1.97	0.46
1:D:36:GLN:NE2	2:D:242:E64:O2	2.48	0.46
1:D:158:PHE:O	1:D:159:TYR:C	2.53	0.46
1:D:178:LEU:HA	1:D:202:ILE:HG22	1.98	0.46
1:A:23:ASP:CB	1:A:26:LEU:HD12	2.46	0.45
1:A:132:PRO:HG2	1:A:135:LYS:HB2	1.98	0.45
1:A:49:GLY:HA3	1:A:177:MET:SD	2.56	0.45
1:B:177:MET:HE3	1:B:203:LYS:CD	2.46	0.45
1:C:48:ILE:HD12	1:C:67:GLU:HA	1.98	0.45
1:A:152:VAL:HG22	1:A:156:PHE:CD2	2.51	0.45
1:D:46:SER:CB	1:D:175:ALA:HB1	2.47	0.45
1:B:83:GLY:O	1:B:84:LEU:HD23	2.17	0.45
1:D:65:LEU:HA	1:D:100:PRO:HA	1.99	0.45
1:A:68:GLN:O	1:A:71:VAL:N	2.50	0.44
1:B:148:ILE:HG22	1:B:235:ALA:CB	2.47	0.44
1:A:15:GLU:HB3	1:A:16:ASN:C	2.37	0.44
1:B:152:VAL:HG22	1:B:173:ASN:OD1	2.17	0.44
1:B:51:VAL:CG2	1:B:65:LEU:HD12	2.47	0.44
1:A:229:CYS:O	1:A:231:LEU:N	2.51	0.44
1:A:230:GLY:O	1:A:231:LEU:C	2.55	0.44
1:A:66:SER:HB2	1:A:104:TYR:HB3	1.99	0.44
1:A:125:ILE:HA	1:A:240:ILE:HG22	2.00	0.43
1:A:38:ASN:O	1:A:39:CYS:HB2	2.18	0.43
1:A:36:GLN:NE2	2:A:242:E64:O2	2.51	0.43
1:D:153:SER:O	1:D:155:ASP:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:LEU:HD21	1:B:180:GLY:N	2.34	0.43
1:C:187:VAL:HA	1:C:194:GLY:HA2	2.02	0.42
1:A:52:GLU:OE2	1:A:64:THR:HG22	2.20	0.42
1:D:42:CYS:HB2	1:D:175:ALA:HB3	2.01	0.42
1:B:45:PHE:CE1	1:B:67:GLU:HG3	2.54	0.42
1:B:209:GLN:OE1	1:B:209:GLN:N	2.52	0.42
1:B:181:PHE:CA	1:B:201:ILE:HD12	2.50	0.42
1:B:65:LEU:HD22	1:B:98:ILE:O	2.19	0.42
1:C:209:GLN:H	1:C:209:GLN:CD	2.21	0.42
1:D:232:GLY:O	1:D:233:THR:C	2.59	0.41
1:C:62:LEU:O	1:C:63:ILE:HD13	2.21	0.41
1:C:61:LYS:HZ2	1:C:63:ILE:HD11	1.84	0.41
1:B:89:PHE:O	1:B:93:ILE:HD12	2.20	0.41
1:D:74:SER:OG	1:D:77:ASN:ND2	2.53	0.41
1:D:73:CYS:HB3	1:D:95:LEU:HD21	2.01	0.41
1:A:52:GLU:HG2	1:A:63:ILE:O	2.20	0.41
1:C:65:LEU:HA	1:C:100:PRO:HA	2.02	0.41
1:A:45:PHE:CZ	1:A:67:GLU:HG3	2.55	0.41
1:C:139:ALA:HB3	1:C:146:ILE:HD11	2.02	0.41
1:C:139:ALA:HB1	1:C:146:ILE:HD11	2.03	0.41
1:C:149:SER:OG	2:C:242:E64:H103	2.21	0.41
1:C:85:ILE:O	1:C:88:ALA:HB3	2.21	0.41
1:C:62:LEU:C	1:C:62:LEU:HD23	2.41	0.40
1:B:230:GLY:O	1:B:232:GLY:N	2.53	0.40
1:A:146:ILE:O	1:A:148:ILE:HG23	2.21	0.40
1:A:230:GLY:O	1:A:232:GLY:N	2.53	0.40
1:A:100:PRO:O	1:A:102:GLY:N	2.55	0.40
1:A:38:ASN:C	1:A:38:ASN:OD1	2.59	0.40
1:A:52:GLU:OE2	1:A:64:THR:HA	2.20	0.40
1:B:31:THR:HG21	1:B:52:GLU:HB3	2.02	0.40
1:B:36:GLN:HG3	1:B:37:LYS:N	2.36	0.40
1:B:93:ILE:HD12	1:B:125:ILE:HD11	2.04	0.40
1:A:178:LEU:HD23	1:A:202:ILE:HD11	2.00	0.40
1:A:65:LEU:HA	1:A:100:PRO:HA	2.03	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	232/241 (96%)	205 (88%)	20 (9%)	7 (3%)	5	22
1	B	238/241 (99%)	217 (91%)	18 (8%)	3 (1%)	15	46
1	C	230/241 (95%)	198 (86%)	28 (12%)	4 (2%)	11	38
1	D	209/241 (87%)	172 (82%)	29 (14%)	8 (4%)	4	16
All	All	909/964 (94%)	792 (87%)	95 (10%)	22 (2%)	7	29

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	39	CYS
1	C	163	ILE
1	C	192	LYS
1	D	154	ASP
1	A	101	ASP
1	A	213	ARG
1	A	230	GLY
1	B	231	LEU
1	C	231	LEU
1	D	40	GLY
1	D	208	GLN
1	D	231	LEU
1	A	61	LYS
1	A	231	LEU
1	B	37	LYS
1	B	190	LEU
1	C	160	LYS
1	D	39	CYS
1	D	159	TYR
1	D	165	ASP
1	D	134	ASN
1	A	150	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	201/206 (98%)	188 (94%)	13 (6%)	21	52
1	B	205/206 (100%)	197 (96%)	8 (4%)	39	75
1	C	200/206 (97%)	190 (95%)	10 (5%)	30	65
1	D	183/206 (89%)	172 (94%)	11 (6%)	24	57
All	All	789/824 (96%)	747 (95%)	42 (5%)	28	63

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	GLN
1	A	3	ASN
1	A	33	VAL
1	A	38	ASN
1	A	51	VAL
1	A	64	THR
1	A	104	TYR
1	A	107	VAL
1	A	155	ASP
1	A	183	MET
1	A	208	GLN
1	A	233	THR
1	A	237	ILE
1	B	51	VAL
1	B	62	LEU
1	B	63	ILE
1	B	67	GLU
1	B	68	GLN
1	B	104	TYR
1	B	183	MET
1	B	233	THR
1	C	64	THR
1	C	90	GLU
1	C	104	TYR

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Mol	Chain	Res	Type
1	C	108	SER
1	C	112	ASN
1	C	120	THR
1	C	125	ILE
1	C	171	GLN
1	C	209	GLN
1	C	233	THR
1	D	31	THR
1	D	55	TYR
1	D	90	GLU
1	D	103	ASP
1	D	104	TYR
1	D	120	THR
1	D	129	LEU
1	D	154	ASP
1	D	181	PHE
1	D	216	ILE
1	D	241	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	68	GLN
1	C	38	ASN
1	D	112	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

5 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	E64	A	242	1	17,24,24	0.43	0	20,30,30	0.81	0
2	E64	B	242	1	17,24,24	0.42	0	20,30,30	0.71	0
2	E64	C	242	1	17,24,24	0.55	0	20,30,30	0.80	1 (5%)
3	GOL	C	243	-	5,5,5	0.35	0	5,5,5	0.22	0
2	E64	D	242	1	17,24,24	0.47	0	20,30,30	0.61	0

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	E64	A	242	1	-	0/25/29/29	0/0/0/0
2	E64	B	242	1	-	0/25/29/29	0/0/0/0
2	E64	C	242	1	-	0/25/29/29	0/0/0/0
3	GOL	C	243	-	-	0/4/4/4	0/0/0/0
2	E64	D	242	1	-	0/25/29/29	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	242	E64	C2-C3-C4	-2.23	106.91	111.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	242	E64	1	0
2	C	242	E64	1	0
2	D	242	E64	1	0

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	236/241 (97%)	0.62	10 (4%) 40 33	61, 73, 80, 85	24 (10%)
1	B	240/241 (99%)	0.48	12 (5%) 32 26	73, 79, 90, 94	47 (19%)
1	C	234/241 (97%)	0.58	14 (5%) 25 18	68, 74, 78, 87	52 (22%)
1	D	215/241 (89%)	0.24	11 (5%) 32 25	58, 83, 101, 115	66 (30%)
All	All	925/964 (95%)	0.49	47 (5%) 32 25	58, 75, 92, 115	189 (20%)

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	142	PHE	4.9
1	D	168	CYS	4.5
1	D	179	VAL	4.1
1	C	200	TYR	3.3
1	B	210	TRP	3.1
1	B	204	ASN	3.1
1	B	175	ALA	3.0
1	C	178	LEU	2.9
1	A	147	SER	2.9
1	D	20	ALA	2.9
1	C	165	ASP	2.8
1	B	150	VAL	2.8
1	C	169	GLY	2.7
1	C	152	VAL	2.7
1	A	70	LEU	2.6
1	A	98	ILE	2.6
1	C	143	LEU	2.5
1	D	167	GLU	2.5
1	C	183	MET	2.5
1	B	75	PHE	2.5
1	B	70	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	116	ILE	2.4
1	D	231	LEU	2.4
1	A	46	SER	2.4
1	D	214	GLY	2.4
1	B	142	PHE	2.4
1	C	202	ILE	2.3
1	C	104	TYR	2.3
1	C	164	PHE	2.3
1	B	231	LEU	2.3
1	D	143	LEU	2.3
1	A	146	ILE	2.3
1	D	180	GLY	2.3
1	A	149	SER	2.2
1	C	118	ARG	2.2
1	B	107	VAL	2.2
1	B	41	SER	2.2
1	A	201	ILE	2.2
1	D	150	VAL	2.1
1	A	63	ILE	2.1
1	C	238	PRO	2.1
1	D	157	ALA	2.1
1	D	51	VAL	2.1
1	B	214	GLY	2.1
1	A	218	ILE	2.1
1	B	66	SER	2.1
1	A	89	PHE	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 carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	E64	B	242	25/25	0.84	0.40	1.94	76,79,80,81	7
2	E64	C	242	25/25	0.87	0.29	1.17	79,83,87,87	0
2	E64	D	242	25/25	0.89	0.19	0.30	69,72,78,79	0
2	E64	A	242	25/25	0.84	0.27	0.21	69,72,80,80	0
3	GOL	C	243	6/6	0.64	0.19	-0.00	123,124,124,124	0

6.5 Other polymers [i](#)

There are no such residues in this entry.