



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 22, 2018 – 04:22 PM EST

PDB ID : 5OSB
Title : GLIC-GABAAR alpha1 chimera crystallized in complex with THDOC at pH4.5
Authors : Lavery, D.C.; Gold, M.G.; Smart, T.G.
Deposited on : 2017-08-17
Resolution : 3.80 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

<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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20030736

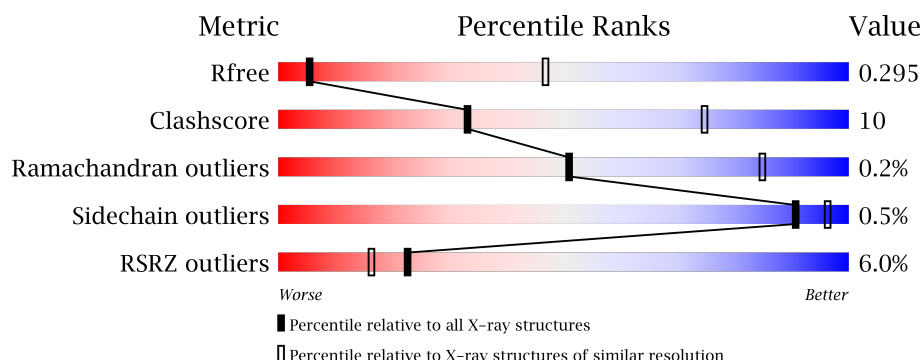
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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}	100719	1019 (4.08-3.52)
Clashscore	112137	1030 (4.04-3.56)
Ramachandran outliers	110173	1011 (4.06-3.54)
Sidechain outliers	110143	1005 (4.06-3.54)
RSRZ outliers	101464	1032 (4.08-3.52)

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	336	<div> <div>2%</div> <div>75% 18% 7%</div> </div>
1	B	336	<div> <div>5%</div> <div>70% 23% 7%</div> </div>
1	C	336	<div> <div>12%</div> <div>72% 20% 8%</div> </div>
1	D	336	<div> <div>5%</div> <div>73% 20% 7%</div> </div>
1	E	336	<div> <div>4%</div> <div>73% 20% 7%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	A8Z	A	501	-	-	-	X
2	A8Z	B	501	-	-	-	X
2	A8Z	C	501	-	-	-	X
2	A8Z	E	501	-	-	-	X
3	ACT	B	502	-	-	X	X
3	ACT	C	502	-	-	-	X
3	ACT	D	501	-	-	-	X
3	ACT	E	502	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12746 atoms, of which 185 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 Proton-gated ion channel, Gamma-aminobutyric acid receptor subunit alpha-1, Gamma-aminobutyric acid receptor subunit alpha-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2485	1624	399	456	6			
1	B	313	Total	C	N	O	S	0	0	0
			2491	1627	402	456	6			
1	C	310	Total	C	N	O	S	0	0	0
			2473	1617	397	453	6			
1	D	313	Total	C	N	O	S	0	0	0
			2485	1624	399	456	6			
1	E	313	Total	C	N	O	S	0	0	0
			2485	1624	399	456	6			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	VAL	GLY	conflict	UNP P62812
A	312	SER	-	linker	UNP P62812
A	313	GLN	-	linker	UNP P62812
A	314	PRO	-	linker	UNP P62812
A	315	ALA	-	linker	UNP P62812
A	316	ARG	-	linker	UNP P62812
A	317	ALA	-	linker	UNP P62812
A	318	ALA	-	linker	UNP P62812
A	429	HIS	-	expression tag	UNP P62812
A	430	HIS	-	expression tag	UNP P62812
A	431	HIS	-	expression tag	UNP P62812
A	432	HIS	-	expression tag	UNP P62812
A	433	HIS	-	expression tag	UNP P62812
A	434	HIS	-	expression tag	UNP P62812
A	435	HIS	-	expression tag	UNP P62812
A	436	HIS	-	expression tag	UNP P62812
B	258	VAL	GLY	conflict	UNP P62812
B	312	SER	-	linker	UNP P62812

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Chain	Residue	Modelled	Actual	Comment	Reference
B	313	GLN	-	linker	UNP P62812
B	314	PRO	-	linker	UNP P62812
B	315	ALA	-	linker	UNP P62812
B	316	ARG	-	linker	UNP P62812
B	317	ALA	-	linker	UNP P62812
B	318	ALA	-	linker	UNP P62812
B	429	HIS	-	expression tag	UNP P62812
B	430	HIS	-	expression tag	UNP P62812
B	431	HIS	-	expression tag	UNP P62812
B	432	HIS	-	expression tag	UNP P62812
B	433	HIS	-	expression tag	UNP P62812
B	434	HIS	-	expression tag	UNP P62812
B	435	HIS	-	expression tag	UNP P62812
B	436	HIS	-	expression tag	UNP P62812
C	258	VAL	GLY	conflict	UNP P62812
C	312	SER	-	linker	UNP P62812
C	313	GLN	-	linker	UNP P62812
C	314	PRO	-	linker	UNP P62812
C	315	ALA	-	linker	UNP P62812
C	316	ARG	-	linker	UNP P62812
C	317	ALA	-	linker	UNP P62812
C	318	ALA	-	linker	UNP P62812
C	429	HIS	-	expression tag	UNP P62812
C	430	HIS	-	expression tag	UNP P62812
C	431	HIS	-	expression tag	UNP P62812
C	432	HIS	-	expression tag	UNP P62812
C	433	HIS	-	expression tag	UNP P62812
C	434	HIS	-	expression tag	UNP P62812
C	435	HIS	-	expression tag	UNP P62812
C	436	HIS	-	expression tag	UNP P62812
D	258	VAL	GLY	conflict	UNP P62812
D	312	SER	-	linker	UNP P62812
D	313	GLN	-	linker	UNP P62812
D	314	PRO	-	linker	UNP P62812
D	315	ALA	-	linker	UNP P62812
D	316	ARG	-	linker	UNP P62812
D	317	ALA	-	linker	UNP P62812
D	318	ALA	-	linker	UNP P62812
D	429	HIS	-	expression tag	UNP P62812
D	430	HIS	-	expression tag	UNP P62812
D	431	HIS	-	expression tag	UNP P62812
D	432	HIS	-	expression tag	UNP P62812

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Chain	Residue	Modelled	Actual	Comment	Reference
D	433	HIS	-	expression tag	UNP P62812
D	434	HIS	-	expression tag	UNP P62812
D	435	HIS	-	expression tag	UNP P62812
D	436	HIS	-	expression tag	UNP P62812
E	258	VAL	GLY	conflict	UNP P62812
E	312	SER	-	linker	UNP P62812
E	313	GLN	-	linker	UNP P62812
E	314	PRO	-	linker	UNP P62812
E	315	ALA	-	linker	UNP P62812
E	316	ARG	-	linker	UNP P62812
E	317	ALA	-	linker	UNP P62812
E	318	ALA	-	linker	UNP P62812
E	429	HIS	-	expression tag	UNP P62812
E	430	HIS	-	expression tag	UNP P62812
E	431	HIS	-	expression tag	UNP P62812
E	432	HIS	-	expression tag	UNP P62812
E	433	HIS	-	expression tag	UNP P62812
E	434	HIS	-	expression tag	UNP P62812
E	435	HIS	-	expression tag	UNP P62812
E	436	HIS	-	expression tag	UNP P62812

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- The chemical structure of A8Z is a complex polycyclic molecule. It features a central core with several fused and fused rings. Key features include:
- Functional Groups:** A hydroxyl group (OH) is attached to C1, and a carboxylic acid group (COOH) is attached to C17. Both are highlighted in red.
 - Stereocenters:** Numerous chiral centers are indicated by wedged and dashed bonds. For example, C10(S), C13(S), C14(S), C17(S), C19, C20, C21, C2, C3(R), C4, C5(S), C6, C7, C8(R), C9(S), C11, C12, C15, C16, C18, and C19 are labeled with their stereochemistry.
 - Hydrogens:** Several hydrogens are explicitly shown, including H1, H2, H3, H4, and H5, with their stereochemistry indicated by wedged or dashed bonds.
 - Carbons:** The carbon skeleton is extensively labeled with C1 through C21, indicating the positions of various atoms and functional groups.

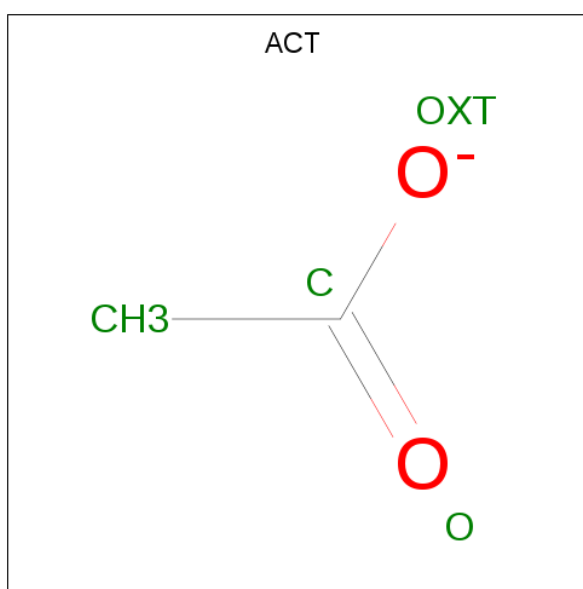
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			58	21	34	3		



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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			58	21	34	3		
2	B	1	Total	C	H	O	0	0
			58	21	34	3		
2	C	1	Total	C	H	O	0	0
			58	21	34	3		
2	E	1	Total	C	H	O	0	0
			58	21	34	3		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).

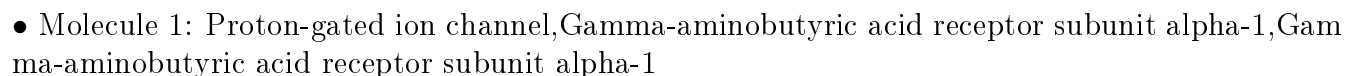
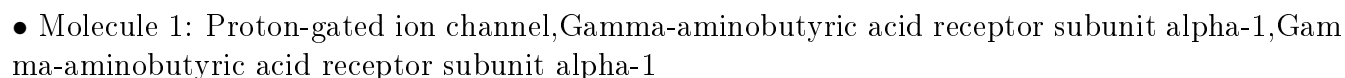


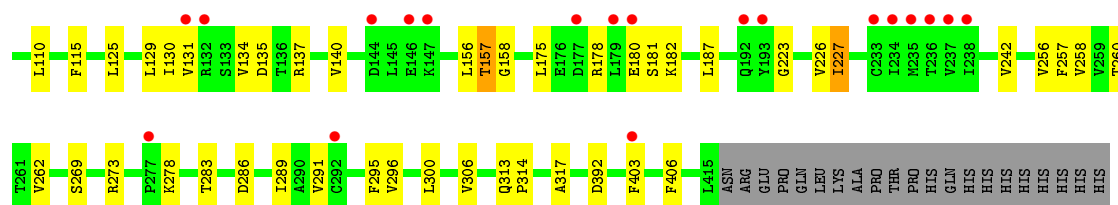
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			7	2	3	2		
3	B	1	Total	C	H	O	0	0
			7	2	3	2		
3	C	1	Total	C	H	O	0	0
			7	2	3	2		
3	D	1	Total	C	H	O	0	0
			7	2	3	2		
3	E	1	Total	C	H	O	0	0
			7	2	3	2		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

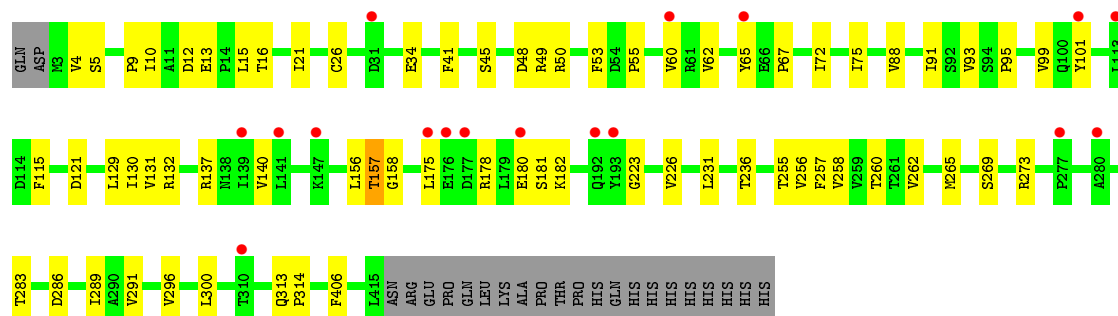
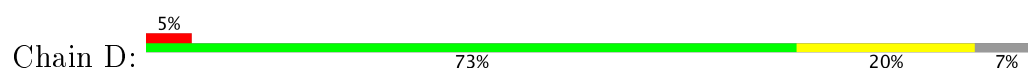
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Cl 1	0	0
4	D	1	Total 1	Cl 1	0	0

- Molecule 1: Proton-gated ion channel, Gamma-aminobutyric acid receptor subunit alpha-1, Gamma-aminobutyric acid receptor subunit alpha-1

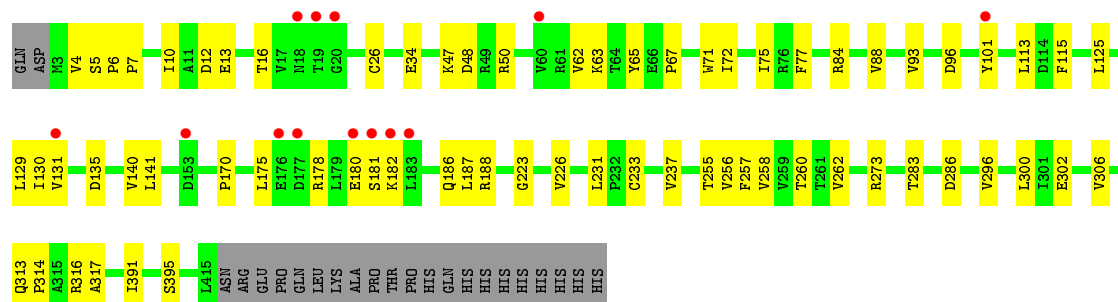
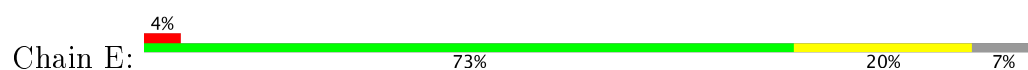




- Molecule 1: Proton-gated ion channel, Gamma-aminobutyric acid receptor subunit alpha-1, Gamma-aminobutyric acid receptor subunit alpha-1



- Molecule 1: Proton-gated ion channel, Gamma-aminobutyric acid receptor subunit alpha-1, Gamma-aminobutyric acid receptor subunit alpha-1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.22Å 133.51Å 162.26Å 90.00° 103.41° 90.00°	Depositor
Resolution (Å)	29.98 – 3.80 87.46 – 3.80	Depositor EDS
% Data completeness (in resolution range)	97.2 (29.98-3.80) 99.4 (87.46-3.80)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 3.78Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.241 , 0.290 0.240 , 0.295	Depositor DCC
R_{free} test set	1901 reflections (5.14%)	DCC
Wilson B-factor (Å ²)	91.0	Xtriage
Anisotropy	0.771	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 99.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	12746	wwPDB-VP
Average B, all atoms (Å ²)	189.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.41% 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: A8Z, CL, ACT

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	0.24	0/2550	0.40	0/3495
1	B	0.24	0/2556	0.40	0/3502
1	C	0.24	0/2537	0.40	0/3474
1	D	0.24	0/2550	0.40	0/3495
1	E	0.24	0/2550	0.40	0/3495
All	All	0.24	0/12743	0.40	0/17461

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	2485	0	2443	39	0
1	B	2491	0	2454	60	0
1	C	2473	0	2439	57	0
1	D	2485	0	2443	51	0
1	E	2485	0	2443	52	0
2	A	48	68	0	0	0
2	B	24	34	0	2	0
2	C	24	34	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	24	34	0	0	0
3	A	4	3	3	1	0
3	B	4	3	3	2	0
3	C	4	3	3	1	0
3	D	4	3	3	1	0
3	E	4	3	3	1	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
All	All	12561	185	12237	244	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 10.

All (244) 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:B:76:ARG:NH1	3:B:502:ACT:O	2.20	0.74
1:B:10:ILE:HB	1:B:13:GLU:HB2	1.71	0.72
1:B:227:ILE:HD11	1:C:273:ARG:CZ	2.20	0.71
1:D:48:ASP:OD1	1:D:50:ARG:HD3	1.90	0.70
1:D:88:VAL:HG22	1:D:101:TYR:OH	1.91	0.70
1:D:9:PRO:HG3	1:D:49:ARG:CZ	2.23	0.69
1:D:130:ILE:HD13	3:D:501:ACT:H2	1.76	0.68
1:D:156:LEU:O	1:D:158:GLY:N	2.27	0.67
1:D:55:PRO:HG3	1:D:95:PRO:HB3	1.77	0.67
1:C:156:LEU:O	1:C:158:GLY:N	2.29	0.65
1:E:10:ILE:HB	1:E:13:GLU:HB2	1.79	0.64
1:B:88:VAL:HG22	1:B:101:TYR:OH	1.98	0.64
1:B:256:VAL:O	1:B:260:THR:HG23	1.98	0.63
1:B:62:VAL:HG23	1:B:63:LYS:N	2.15	0.62
1:B:10:ILE:O	1:B:10:ILE:HG22	1.99	0.62
1:E:129:LEU:O	1:E:182:LYS:HA	1.99	0.61
1:B:306:VAL:HG13	1:B:317:ALA:HB1	1.82	0.61
1:E:62:VAL:HG23	1:E:63:LYS:N	2.16	0.61
1:B:130:ILE:HD13	3:B:502:ACT:H2	1.83	0.60
1:C:227:ILE:HD11	1:D:273:ARG:CZ	2.32	0.59
1:D:157:THR:HG21	1:E:34:GLU:OE1	2.03	0.59
1:B:48:ASP:OD1	1:B:50:ARG:HD3	2.02	0.59
1:E:48:ASP:OD1	1:E:50:ARG:HD3	2.02	0.59
1:C:16:THR:HA	1:C:140:VAL:O	2.04	0.58
1:B:65:TYR:O	1:B:91:ILE:HB	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:75:ILE:HD13	1:D:131:VAL:HB	1.86	0.58
1:D:130:ILE:HA	1:D:181:SER:O	2.05	0.57
1:D:55:PRO:HG3	1:D:95:PRO:CB	2.34	0.57
1:E:258:VAL:HG11	1:E:300:LEU:CD1	2.35	0.57
1:C:157:THR:HG21	1:D:34:GLU:OE1	2.04	0.57
1:C:258:VAL:HG11	1:C:300:LEU:CD1	2.35	0.57
1:C:7:PRO:HD2	1:C:50:ARG:HD2	1.87	0.57
1:E:88:VAL:HG22	1:E:101:TYR:OH	2.05	0.57
1:C:75:ILE:HD13	1:C:131:VAL:HB	1.87	0.56
1:D:62:VAL:HG13	1:D:93:VAL:O	2.05	0.56
1:D:16:THR:HA	1:D:140:VAL:O	2.05	0.56
1:B:224:TYR:CD2	1:C:278:LYS:HA	2.41	0.56
1:E:175:LEU:HB3	1:E:180:GLU:HG3	1.87	0.55
1:E:75:ILE:HD13	1:E:131:VAL:HB	1.89	0.55
1:B:129:LEU:O	1:B:182:LYS:HA	2.07	0.54
1:A:256:VAL:O	1:A:260:THR:HG23	2.07	0.54
1:B:67:PRO:HG3	1:B:88:VAL:HG11	1.90	0.54
1:E:16:THR:HA	1:E:140:VAL:O	2.07	0.54
1:D:256:VAL:O	1:D:260:THR:HG23	2.08	0.54
1:A:295:PHE:HE2	1:A:403:PHE:HA	1.71	0.54
1:A:75:ILE:HD13	1:A:131:VAL:HB	1.89	0.54
1:B:10:ILE:C	1:B:12:ASP:H	2.11	0.54
1:B:67:PRO:HA	1:B:72:ILE:HD11	1.89	0.54
1:E:10:ILE:HG22	1:E:10:ILE:O	2.08	0.54
1:B:130:ILE:HA	1:B:181:SER:O	2.08	0.54
1:C:256:VAL:O	1:C:260:THR:HG23	2.08	0.53
1:C:262:VAL:HG22	1:C:296:VAL:HG12	1.90	0.53
1:A:48:ASP:OD1	1:A:50:ARG:HD3	2.09	0.53
1:C:269:SER:HB2	1:C:289:ILE:HG21	1.89	0.53
1:A:115:PHE:O	1:A:283:THR:HG22	2.08	0.53
1:B:115:PHE:O	1:B:283:THR:HG22	2.08	0.53
1:B:92:SER:OG	1:C:135:ASP:OD2	2.22	0.53
1:D:129:LEU:O	1:D:182:LYS:HA	2.09	0.52
1:A:10:ILE:HG22	1:A:10:ILE:O	2.09	0.52
1:B:262:VAL:HG22	1:B:296:VAL:HG12	1.91	0.52
1:A:60:VAL:HG12	1:A:62:VAL:H	1.74	0.52
1:C:10:ILE:O	1:C:10:ILE:HG22	2.08	0.52
1:E:262:VAL:HG22	1:E:296:VAL:HG12	1.91	0.52
1:D:262:VAL:HG22	1:D:296:VAL:HG12	1.91	0.52
1:A:10:ILE:HB	1:A:13:GLU:HB2	1.92	0.52
1:E:130:ILE:HD13	3:E:502:ACT:H2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:ARG:NH1	1:B:286:ASP:OD2	2.44	0.51
1:B:241:GLN:OE1	2:B:501:A8Z:O2	2.29	0.51
1:C:156:LEU:C	1:C:158:GLY:H	2.13	0.51
1:C:223:GLY:O	1:C:226:VAL:HB	2.11	0.51
1:D:156:LEU:C	1:D:158:GLY:H	2.14	0.51
1:C:15:LEU:HB2	1:C:137:ARG:CZ	2.41	0.51
1:B:186:GLN:OE1	1:B:188:ARG:NH2	2.44	0.50
1:C:129:LEU:O	1:C:182:LYS:HA	2.11	0.50
1:D:67:PRO:HG3	1:D:91:ILE:HD11	1.92	0.50
1:E:256:VAL:O	1:E:260:THR:HG23	2.11	0.50
1:D:175:LEU:HB3	1:D:180:GLU:HG3	1.92	0.50
1:A:132:ARG:CZ	1:A:178:ARG:HB2	2.40	0.50
1:D:258:VAL:HG11	1:D:300:LEU:HD12	1.93	0.50
1:B:6:PRO:HA	1:B:50:ARG:HH11	1.76	0.50
1:B:224:TYR:CE1	1:C:278:LYS:HG2	2.46	0.50
1:D:65:TYR:HE1	1:D:93:VAL:HG21	1.77	0.49
1:E:258:VAL:HG11	1:E:300:LEU:HD12	1.94	0.49
1:A:15:LEU:HB2	1:A:137:ARG:CZ	2.43	0.49
1:D:10:ILE:C	1:D:12:ASP:H	2.16	0.49
1:D:257:PHE:CE1	1:E:255:THR:HG23	2.48	0.49
1:E:4:VAL:O	1:E:50:ARG:NH2	2.43	0.49
1:D:15:LEU:HB2	1:D:137:ARG:CZ	2.42	0.49
1:E:130:ILE:HA	1:E:181:SER:O	2.11	0.49
1:A:130:ILE:HA	1:A:181:SER:O	2.12	0.49
1:B:16:THR:HA	1:B:140:VAL:O	2.12	0.49
1:C:175:LEU:HB3	1:C:180:GLU:HG3	1.95	0.49
1:E:7:PRO:HD2	1:E:50:ARG:HD2	1.94	0.49
1:B:258:VAL:HG11	1:B:300:LEU:HD12	1.92	0.49
1:C:291:VAL:HG11	1:C:406:PHE:CE1	2.48	0.49
1:E:62:VAL:HG23	1:E:63:LYS:H	1.77	0.49
1:B:37:LYS:HE3	1:B:108:ARG:HD2	1.95	0.49
1:C:313:GLN:N	1:C:314:PRO:HD3	2.28	0.49
1:C:295:PHE:HE2	1:C:403:PHE:HA	1.78	0.49
1:E:6:PRO:HA	1:E:50:ARG:HH11	1.78	0.49
1:A:258:VAL:HG11	1:A:300:LEU:CD1	2.42	0.48
1:E:302:GLU:HG3	1:E:395:SER:OG	2.13	0.48
1:A:75:ILE:CD1	1:A:131:VAL:HB	2.43	0.48
1:A:262:VAL:HG22	1:A:296:VAL:HG12	1.95	0.48
1:B:258:VAL:HG11	1:B:300:LEU:CD1	2.43	0.48
1:D:10:ILE:O	1:D:10:ILE:HG22	2.14	0.48
1:A:65:TYR:HE1	1:A:93:VAL:HG21	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:258:VAL:CG1	1:C:300:LEU:HD13	2.44	0.48
1:C:258:VAL:HG11	1:C:300:LEU:HD13	1.95	0.48
1:B:62:VAL:CG2	1:B:63:LYS:N	2.77	0.48
1:D:291:VAL:HG11	1:D:406:PHE:CE1	2.48	0.48
1:D:157:THR:HG22	1:D:157:THR:O	2.13	0.47
1:E:313:GLN:N	1:E:314:PRO:HD3	2.29	0.47
1:A:291:VAL:HG21	1:A:406:PHE:CE1	2.49	0.47
1:D:132:ARG:CZ	1:D:178:ARG:HB2	2.44	0.47
1:A:129:LEU:O	1:A:182:LYS:HA	2.15	0.47
1:D:60:VAL:HG12	1:D:62:VAL:H	1.80	0.47
1:A:313:GLN:N	1:A:314:PRO:HD3	2.30	0.47
1:E:115:PHE:O	1:E:283:THR:HG22	2.14	0.47
1:B:119:PRO:HD3	1:B:285:MET:HB2	1.96	0.47
1:E:67:PRO:HA	1:E:72:ILE:HD11	1.95	0.47
1:A:306:VAL:HG21	1:A:392:ASP:OD1	2.15	0.47
1:D:53:PHE:CE1	1:D:95:PRO:HA	2.50	0.46
1:A:41:PHE:CE1	1:A:104:ARG:HB2	2.50	0.46
1:E:75:ILE:CD1	1:E:131:VAL:HB	2.45	0.46
1:B:224:TYR:CD1	1:C:278:LYS:HG2	2.51	0.46
1:B:242:VAL:HG22	2:B:501:A8Z:C7	2.45	0.46
1:D:67:PRO:HA	1:D:72:ILE:HD11	1.97	0.46
1:E:4:VAL:HG22	1:E:5:SER:H	1.80	0.46
1:B:295:PHE:HE2	1:B:403:PHE:HA	1.80	0.46
1:A:16:THR:HA	1:A:140:VAL:O	2.16	0.46
1:A:258:VAL:HG11	1:A:300:LEU:HD12	1.96	0.46
1:B:65:TYR:HE2	1:B:93:VAL:HG21	1.81	0.46
1:D:258:VAL:HG11	1:D:300:LEU:CD1	2.46	0.46
1:B:9:PRO:HG3	1:B:49:ARG:HD3	1.97	0.46
1:C:273:ARG:NH1	1:C:286:ASP:OD2	2.49	0.46
1:A:33:ALA:O	1:A:35:THR:HG23	2.16	0.46
1:B:223:GLY:O	1:B:226:VAL:HB	2.16	0.46
1:C:29:LEU:HD13	1:C:36:PHE:HB3	1.98	0.46
1:C:90:ASP:OD2	1:D:178:ARG:NH1	2.36	0.46
1:D:75:ILE:CD1	1:D:131:VAL:HB	2.45	0.46
1:B:7:PRO:CD	1:B:50:ARG:HD2	2.46	0.46
1:C:283:THR:O	1:C:286:ASP:HB2	2.16	0.45
1:E:135:ASP:OD1	1:E:178:ARG:NH2	2.49	0.45
1:B:41:PHE:HE1	1:B:104:ARG:HB2	1.81	0.45
1:D:313:GLN:N	1:D:314:PRO:HD3	2.32	0.45
1:E:10:ILE:C	1:E:12:ASP:H	2.19	0.45
1:A:248:ARG:HD2	1:A:306:VAL:CG1	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ARG:HD2	3:A:503:ACT:O	2.16	0.45
1:C:75:ILE:CD1	1:C:131:VAL:HB	2.45	0.45
1:A:283:THR:O	1:A:286:ASP:HB2	2.16	0.45
1:B:41:PHE:CE1	1:B:104:ARG:HB2	2.51	0.45
1:B:67:PRO:HG3	1:B:91:ILE:HD11	1.98	0.45
1:C:67:PRO:HA	1:C:72:ILE:HD11	1.99	0.45
1:E:7:PRO:CD	1:E:50:ARG:HD2	2.47	0.45
1:C:130:ILE:HA	1:C:181:SER:O	2.17	0.45
1:A:255:THR:HG23	1:E:257:PHE:CE1	2.52	0.45
1:A:47:LYS:HE3	1:A:96:ASP:OD2	2.17	0.45
1:B:135:ASP:OD1	1:B:178:ARG:NH2	2.50	0.45
1:C:10:ILE:HB	1:C:13:GLU:HB2	1.98	0.45
1:D:4:VAL:HG22	1:D:5:SER:H	1.82	0.45
1:E:47:LYS:HE3	1:E:96:ASP:OD2	2.17	0.45
1:E:65:TYR:HE1	1:E:93:VAL:HG21	1.82	0.45
1:A:6:PRO:HA	1:A:50:ARG:NH1	2.32	0.44
1:C:41:PHE:CE1	1:C:104:ARG:HB2	2.52	0.44
1:D:223:GLY:O	1:D:226:VAL:HB	2.17	0.44
1:C:115:PHE:O	1:C:283:THR:HG22	2.18	0.44
1:D:21:ILE:HA	1:D:41:PHE:O	2.17	0.44
1:B:227:ILE:HD11	1:C:273:ARG:NE	2.32	0.44
1:D:283:THR:O	1:D:286:ASP:HB2	2.17	0.44
1:B:75:ILE:HD13	1:B:131:VAL:HB	2.00	0.44
1:C:8:PRO:HG2	1:C:11:ALA:HA	1.99	0.44
1:E:316:ARG:O	1:E:391:ILE:HG13	2.18	0.44
1:C:135:ASP:OD1	1:C:178:ARG:NH2	2.51	0.44
1:C:125:LEU:HD12	1:C:187:LEU:HD23	1.99	0.44
1:D:121:ASP:OD1	1:D:121:ASP:N	2.50	0.43
1:B:25:GLU:HG3	1:C:110:LEU:HD23	2.00	0.43
1:E:231:LEU:HA	1:E:231:LEU:HD23	1.88	0.43
1:E:258:VAL:CG1	1:E:300:LEU:HD13	2.48	0.43
1:B:53:PHE:CE1	1:B:95:PRO:HA	2.53	0.43
1:B:21:ILE:HA	1:B:41:PHE:O	2.18	0.43
1:C:10:ILE:C	1:C:12:ASP:H	2.21	0.43
1:E:129:LEU:HD23	1:E:129:LEU:HA	1.88	0.43
1:A:4:VAL:HG22	1:A:5:SER:H	1.83	0.43
1:C:257:PHE:CE1	1:D:255:THR:HG23	2.54	0.43
1:C:6:PRO:HD3	1:C:71:TRP:CE3	2.54	0.43
1:C:258:VAL:HG13	1:C:300:LEU:HB2	2.01	0.43
1:C:258:VAL:HG11	1:C:300:LEU:HD12	2.01	0.43
1:C:269:SER:HB2	1:C:289:ILE:CG2	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:141:LEU:O	1:E:170:PRO:HG3	2.18	0.43
1:B:269:SER:HB2	1:B:289:ILE:HG21	2.01	0.42
1:C:4:VAL:HG22	1:C:5:SER:H	1.84	0.42
1:E:273:ARG:NH1	1:E:286:ASP:OD2	2.52	0.42
1:C:41:PHE:HE1	1:C:104:ARG:HB2	1.84	0.42
1:D:115:PHE:O	1:D:283:THR:HG22	2.19	0.42
1:E:77:PHE:CE2	1:E:84:ARG:HD3	2.54	0.42
1:A:118:TYR:CD2	1:A:119:PRO:HA	2.54	0.42
1:B:4:VAL:HG22	1:B:5:SER:H	1.84	0.42
1:A:10:ILE:C	1:A:12:ASP:H	2.22	0.42
1:E:6:PRO:HD3	1:E:71:TRP:CE3	2.54	0.42
1:A:6:PRO:HA	1:A:50:ARG:HH11	1.84	0.42
1:A:129:LEU:HA	1:A:129:LEU:HD23	1.88	0.42
1:E:129:LEU:O	1:E:182:LYS:HG3	2.20	0.42
1:A:37:LYS:HE3	1:A:108:ARG:HD2	2.02	0.42
1:A:22:TYR:O	1:A:40:ALA:HB1	2.20	0.42
1:B:283:THR:O	1:B:286:ASP:HB2	2.20	0.42
1:E:306:VAL:HG13	1:E:317:ALA:HB1	2.01	0.42
1:B:7:PRO:HD2	1:B:50:ARG:HD2	2.01	0.42
1:E:62:VAL:CG2	1:E:63:LYS:N	2.81	0.42
1:B:41:PHE:CE2	3:C:502:ACT:H3	2.54	0.42
1:C:242:VAL:HG22	2:C:501:A8Z:C7	2.50	0.42
1:C:306:VAL:HG21	1:C:392:ASP:OD1	2.19	0.42
1:B:313:GLN:N	1:B:314:PRO:HD3	2.34	0.41
1:D:269:SER:HB2	1:D:289:ILE:HG21	2.01	0.41
1:B:291:VAL:HG11	1:B:406:PHE:CE1	2.55	0.41
1:C:157:THR:HG22	1:C:157:THR:O	2.18	0.41
1:D:231:LEU:HA	1:D:231:LEU:HD23	1.90	0.41
1:D:273:ARG:NH1	1:D:286:ASP:OD2	2.53	0.41
1:E:34:GLU:HG2	1:E:113:LEU:HG	2.02	0.41
1:E:258:VAL:HG13	1:E:300:LEU:HB2	2.01	0.41
1:D:236:THR:HG21	1:D:265:MET:HG2	2.01	0.41
1:E:125:LEU:HD12	1:E:187:LEU:HD23	2.02	0.41
1:E:223:GLY:O	1:E:226:VAL:HB	2.19	0.41
1:B:34:GLU:HG2	1:B:113:LEU:HG	2.02	0.41
1:B:54:ASP:O	1:B:58:SER:HB2	2.21	0.41
1:E:283:THR:O	1:E:286:ASP:HB2	2.21	0.41
1:E:6:PRO:HA	1:E:50:ARG:NH1	2.35	0.41
1:A:316:ARG:O	1:A:391:ILE:HG13	2.21	0.41
1:C:129:LEU:HD23	1:C:129:LEU:HA	1.87	0.41
1:D:258:VAL:HG13	1:D:300:LEU:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:186:GLN:OE1	1:E:188:ARG:NH2	2.54	0.41
1:E:233:CYS:O	1:E:237:VAL:HG23	2.21	0.41
1:A:257:PHE:CE1	1:B:255:THR:HG23	2.56	0.41
1:D:45:SER:HA	1:D:99:VAL:O	2.21	0.41
1:C:291:VAL:HG11	1:C:406:PHE:CZ	2.56	0.40
1:D:10:ILE:HB	1:D:13:GLU:HB2	2.03	0.40
1:C:134:VAL:HG23	1:C:137:ARG:H	1.86	0.40
1:B:121:ASP:OD1	1:B:121:ASP:N	2.50	0.40
1:C:306:VAL:HG13	1:C:317:ALA:HB1	2.02	0.40
1:A:295:PHE:CD2	1:A:403:PHE:HD1	2.38	0.40
1:B:125:LEU:HD12	1:B:187:LEU:HD23	2.04	0.40
1:B:4:VAL:O	1:B:50:ARG:NH2	2.51	0.40
1:C:6:PRO:HA	1:C:50:ARG:HH11	1.87	0.40
1:D:55:PRO:HG3	1:D:95:PRO:CG	2.52	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	311/336 (93%)	299 (96%)	12 (4%)	0	100	100
1	B	311/336 (93%)	299 (96%)	12 (4%)	0	100	100
1	C	306/336 (91%)	294 (96%)	10 (3%)	2 (1%)	25	68
1	D	311/336 (93%)	299 (96%)	11 (4%)	1 (0%)	44	81
1	E	311/336 (93%)	299 (96%)	12 (4%)	0	100	100
All	All	1550/1680 (92%)	1490 (96%)	57 (4%)	3 (0%)	51	85

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	157	THR
1	D	157	THR
1	C	63	LYS

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	270/302 (89%)	269 (100%)	1 (0%)	93	97
1	B	271/302 (90%)	269 (99%)	2 (1%)	87	94
1	C	270/302 (89%)	268 (99%)	2 (1%)	87	94
1	D	270/302 (89%)	269 (100%)	1 (0%)	93	97
1	E	270/302 (89%)	269 (100%)	1 (0%)	93	97
All	All	1351/1510 (90%)	1344 (100%)	7 (0%)	91	96

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	CYS
1	B	26	CYS
1	B	227	ILE
1	C	26	CYS
1	C	227	ILE
1	D	26	CYS
1	E	26	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 for Mogul analysis.

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	A8Z	A	501	-	27,27,27	0.26	0	42,43,43	0.63	2 (4%)
2	A8Z	A	502	-	27,27,27	0.26	0	42,43,43	0.59	2 (4%)
3	ACT	A	503	-	1,3,3	1.28	0	0,3,3	0.00	-
2	A8Z	B	501	-	27,27,27	0.25	0	42,43,43	0.63	2 (4%)
3	ACT	B	502	-	1,3,3	1.26	0	0,3,3	0.00	-
2	A8Z	C	501	-	27,27,27	0.25	0	42,43,43	0.63	2 (4%)
3	ACT	C	502	-	1,3,3	1.30	0	0,3,3	0.00	-
3	ACT	D	501	-	1,3,3	1.29	0	0,3,3	0.00	-
2	A8Z	E	501	-	27,27,27	0.26	0	42,43,43	0.60	2 (4%)
3	ACT	E	502	-	1,3,3	1.30	0	0,3,3	0.00	-

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	A8Z	A	501	-	-	0/6/64/64	0/4/4/4
2	A8Z	A	502	-	-	0/6/64/64	0/4/4/4
3	ACT	A	503	-	-	0/0/0/0	0/0/0/0
2	A8Z	B	501	-	-	0/6/64/64	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ACT	B	502	-	-	0/0/0/0	0/0/0/0
2	A8Z	C	501	-	-	0/6/64/64	0/4/4/4
3	ACT	C	502	-	-	0/0/0/0	0/0/0/0
3	ACT	D	501	-	-	0/0/0/0	0/0/0/0
2	A8Z	E	501	-	-	0/6/64/64	0/4/4/4
3	ACT	E	502	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	A8Z	C13-C17-C20	-2.71	111.42	115.46
2	A	501	A8Z	C13-C17-C20	-2.67	111.49	115.46
2	C	501	A8Z	C13-C17-C20	-2.65	111.50	115.46
2	C	501	A8Z	O1-C21-C20	-2.61	105.64	112.66
2	A	501	A8Z	O1-C21-C20	-2.56	105.78	112.66
2	E	501	A8Z	O1-C21-C20	-2.52	105.88	112.66
2	A	502	A8Z	O1-C21-C20	-2.52	105.88	112.66
2	B	501	A8Z	O1-C21-C20	-2.49	105.96	112.66
2	E	501	A8Z	C13-C17-C20	-2.38	111.91	115.46
2	A	502	A8Z	C13-C17-C20	-2.20	112.18	115.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	ACT	1	0
2	B	501	A8Z	2	0
3	B	502	ACT	2	0
2	C	501	A8Z	1	0
3	C	502	ACT	1	0
3	D	501	ACT	1	0
3	E	502	ACT	1	0

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	313/336 (93%)	-0.02	7 (2%) 62 53	130, 180, 227, 331	0
1	B	313/336 (93%)	0.24	17 (5%) 26 21	146, 183, 232, 325	0
1	C	310/336 (92%)	0.53	39 (12%) 4 5	79, 192, 239, 882	0
1	D	313/336 (93%)	0.36	17 (5%) 26 21	140, 183, 238, 300	0
1	E	313/336 (93%)	0.10	13 (4%) 37 29	142, 184, 238, 299	0
All	All	1562/1680 (92%)	0.24	93 (5%) 23 16	79, 184, 237, 882	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	20	GLY	6.0
1	C	147	LYS	5.9
1	D	147	LYS	5.8
1	C	103	GLU	5.8
1	C	19	THR	5.6
1	C	11	ALA	5.5
1	C	43	SER	5.4
1	D	176	GLU	5.4
1	B	147	LYS	5.2
1	E	177	ASP	4.7
1	C	12	ASP	4.7
1	C	42	LEU	4.6
1	C	177	ASP	4.4
1	A	90	ASP	4.4
1	B	397	ILE	4.3
1	A	101	TYR	4.2
1	E	176	GLU	4.2
1	E	19	THR	4.1
1	C	10	ILE	4.0
1	C	144	ASP	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	101	TYR	3.8
1	B	184	ASP	3.7
1	C	132	ARG	3.7
1	D	177	ASP	3.4
1	D	180	GLU	3.3
1	B	60	VAL	3.1
1	E	60	VAL	3.1
1	C	235	MET	3.1
1	C	234	ILE	3.1
1	B	186	GLN	3.1
1	E	183	LEU	3.0
1	D	280	ALA	3.0
1	C	87	ASP	3.0
1	B	398	ALA	2.9
1	C	41	PHE	2.9
1	D	175	LEU	2.9
1	C	104	ARG	2.9
1	D	101	TYR	2.9
1	C	146	GLU	2.8
1	D	65	TYR	2.8
1	C	292	CYS	2.8
1	B	176	GLU	2.8
1	C	237	VAL	2.8
1	B	148	VAL	2.7
1	B	188	ARG	2.7
1	D	310	THR	2.7
1	D	139	ILE	2.6
1	B	394	LEU	2.6
1	B	401	LEU	2.6
1	D	192	GLN	2.6
1	C	193	TYR	2.6
1	C	44	LEU	2.6
1	D	141	LEU	2.5
1	D	60	VAL	2.5
1	C	192	GLN	2.5
1	D	113	LEU	2.5
1	C	180	GLU	2.5
1	E	180	GLU	2.5
1	B	187	LEU	2.5
1	C	236	THR	2.4
1	C	403	PHE	2.4
1	C	179	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	193	TYR	2.4
1	C	18	ASN	2.4
1	C	131	VAL	2.4
1	C	238	ILE	2.3
1	D	31	ASP	2.3
1	A	188	ARG	2.3
1	B	185	TYR	2.3
1	E	182	LYS	2.3
1	A	12	ASP	2.3
1	A	103	GLU	2.3
1	C	86	ALA	2.3
1	E	181	SER	2.3
1	C	233	CYS	2.2
1	B	87	ASP	2.2
1	B	165	THR	2.2
1	D	277	PRO	2.2
1	C	75	ILE	2.2
1	C	102	LEU	2.2
1	E	101	TYR	2.1
1	C	17	VAL	2.1
1	E	153	ASP	2.1
1	C	76	ARG	2.1
1	E	18	ASN	2.1
1	E	20	GLY	2.1
1	B	246	LEU	2.1
1	C	84	ARG	2.1
1	A	100	GLN	2.1
1	A	44	LEU	2.1
1	B	126	HIS	2.0
1	E	131	VAL	2.0
1	C	277	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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(Å ²)	Q<0.9
3	ACT	B	502	4/4	0.93	0.83	6.78	164,167,212,212	0
3	ACT	E	502	4/4	0.92	1.08	5.71	133,156,256,256	0
3	ACT	D	501	4/4	0.86	1.34	4.86	168,202,293,293	0
3	ACT	C	502	4/4	0.90	1.10	3.75	166,184,260,260	0
2	A8Z	A	501	24/24	0.89	0.40	2.58	166,198,254,274	0
3	ACT	A	503	4/4	0.93	0.34	1.95	156,177,188,193	0
2	A8Z	B	501	24/24	0.86	0.54	1.94	176,215,284,291	0
2	A8Z	E	501	24/24	0.89	0.45	1.13	171,207,234,245	0
2	A8Z	C	501	24/24	0.85	0.46	1.06	175,218,272,291	0
2	A8Z	A	502	24/24	0.89	0.24	-0.56	171,214,261,282	0
4	CL	D	502	1/1	0.93	0.22	-1.49	144,144,144,144	0
4	CL	A	504	1/1	0.88	0.18	-2.79	144,144,144,144	0

6.5 Other polymers

There are no such residues in this entry.