



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:32 AM UTC

PDB ID : 2V41 / pdb_00002v41
Title : Crystal Structure of the C45S mutant of the Peroxiredoxin 6 of Arenicola Marina. Orthorhombic form
Authors : Smeets, A.; Declercq, J.P.
Deposited on : 2007-06-27
Resolution : 2.40 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

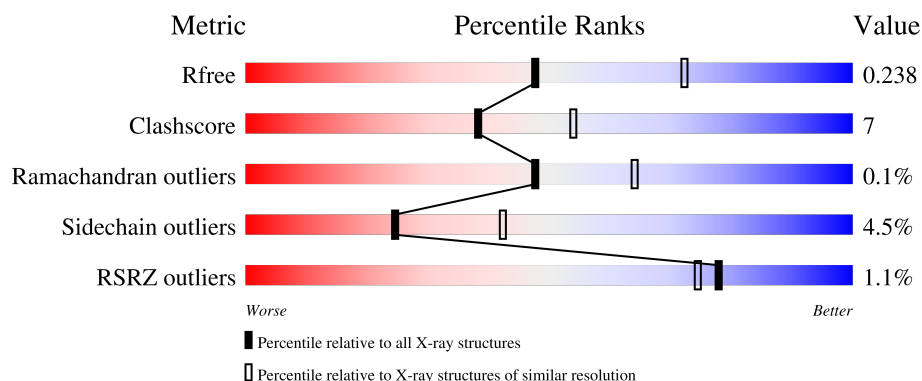
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	233	2% 76% 16% 6%
1	B	233	85% 9% 6%
1	C	233	78% 12% 7%
1	D	233	79% 14% 6%
1	E	233	2% 73% 18% 6%

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Mol	Chain	Length	Quality of chain
1	F	233	2% 80% 13% 6%
1	G	233	2% 75% 18% 6%
1	H	233	% 79% 14% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13959 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.

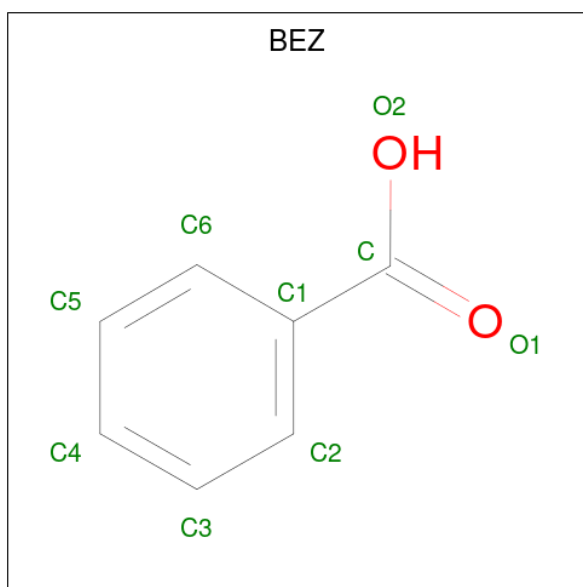
- Molecule 1 is a protein called PEROXIREDOXIN 6..

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	0	0
			1703	1087	288	319	9			
1	B	220	Total	C	N	O	S	0	0	0
			1703	1087	288	319	9			
1	C	216	Total	C	N	O	S	0	0	0
			1672	1068	282	313	9			
1	D	220	Total	C	N	O	S	0	0	0
			1703	1087	288	319	9			
1	E	220	Total	C	N	O	S	0	0	0
			1703	1087	288	319	9			
1	F	220	Total	C	N	O	S	0	0	0
			1703	1087	288	319	9			
1	G	220	Total	C	N	O	S	0	0	0
			1703	1087	288	319	9			
1	H	220	Total	C	N	O	S	0	0	0
			1703	1087	288	319	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	SER	CYS	engineered mutation	UNP Q1AN22
B	45	SER	CYS	engineered mutation	UNP Q1AN22
C	45	SER	CYS	engineered mutation	UNP Q1AN22
D	45	SER	CYS	engineered mutation	UNP Q1AN22
E	45	SER	CYS	engineered mutation	UNP Q1AN22
F	45	SER	CYS	engineered mutation	UNP Q1AN22
G	45	SER	CYS	engineered mutation	UNP Q1AN22
H	45	SER	CYS	engineered mutation	UNP Q1AN22

- Molecule 2 is BENZOIC ACID (CCD ID: BEZ) (formula: C₇H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 9 7 2	0	0
2	B	1	Total C O 9 7 2	0	0
2	C	1	Total C O 9 7 2	0	0
2	D	1	Total C O 9 7 2	0	0
2	E	1	Total C O 9 7 2	0	0
2	F	1	Total C O 9 7 2	0	0
2	G	1	Total C O 9 7 2	0	0
2	H	1	Total C O 9 7 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	46	Total O 46 46	0	0
3	B	48	Total O 48 48	0	0
3	C	33	Total O 33 33	0	0
3	D	39	Total O 39 39	0	0

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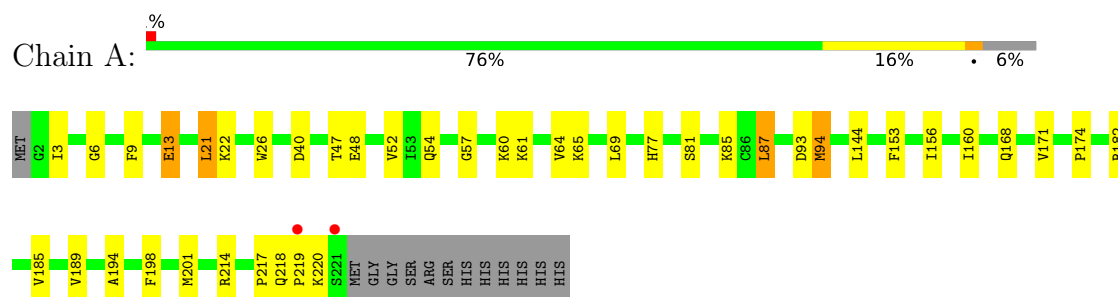
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	31	Total 31	O 31	0	0
3	F	30	Total 30	O 30	0	0
3	G	33	Total 33	O 33	0	0
3	H	34	Total 34	O 34	0	0

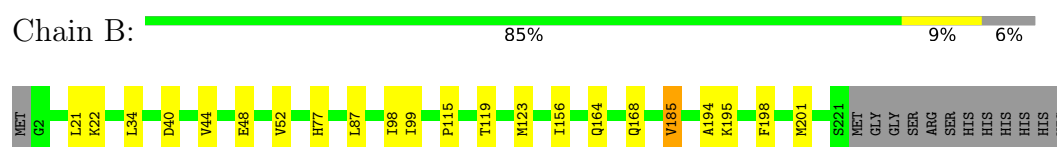
3 Residue-property plots

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.

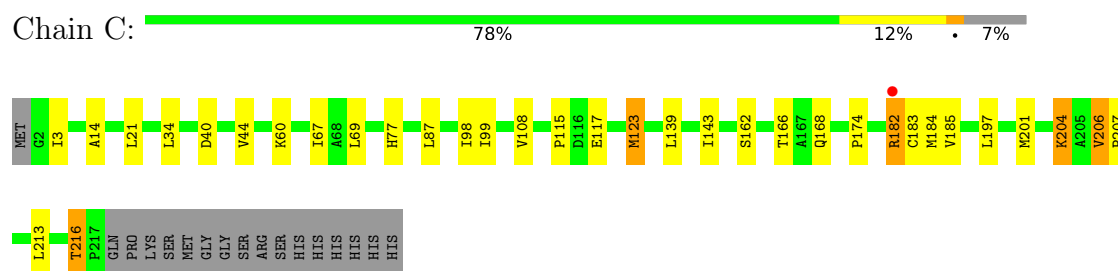
- Molecule 1: PEROXIREDOXIN 6.



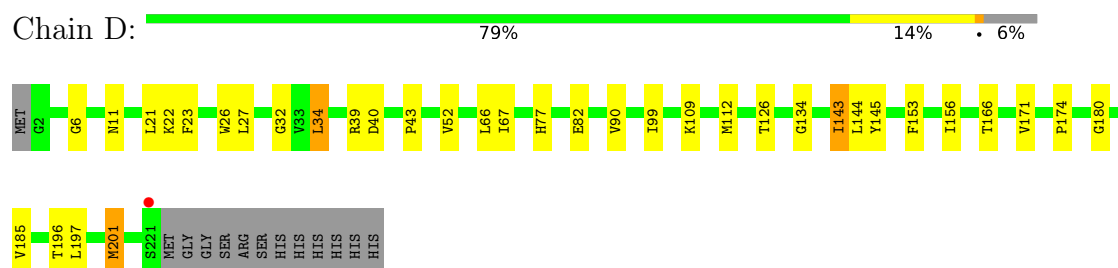
- Molecule 1: PEROXIREDOXIN 6.



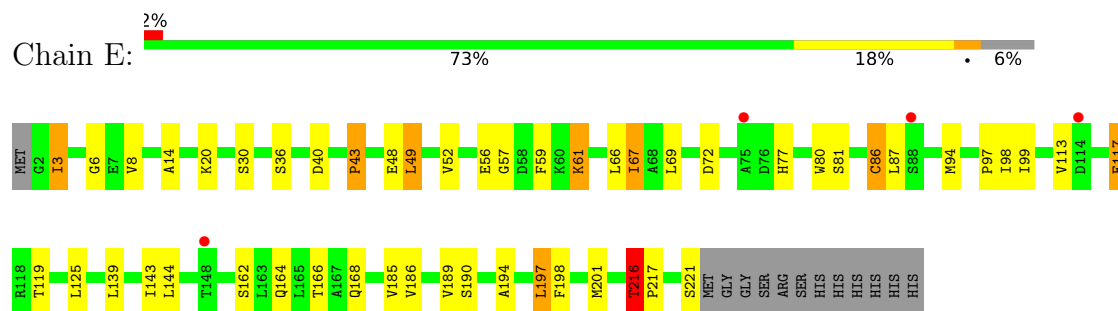
- Molecule 1: PEROXIREDOXIN 6.



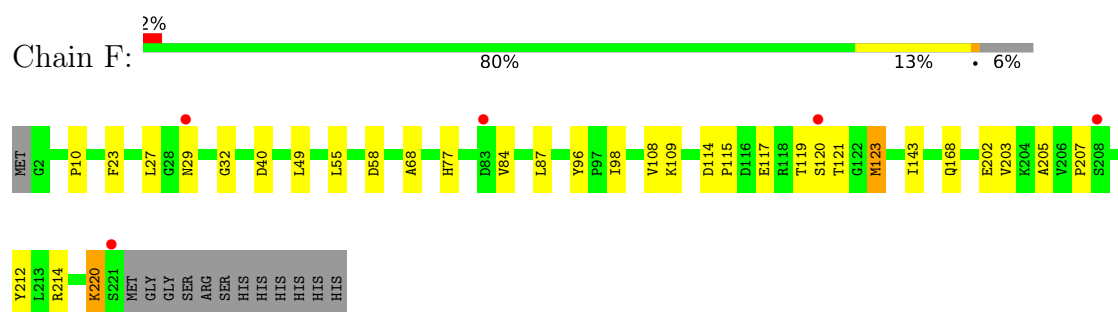
- Molecule 1: PEROXIREDOXIN 6.



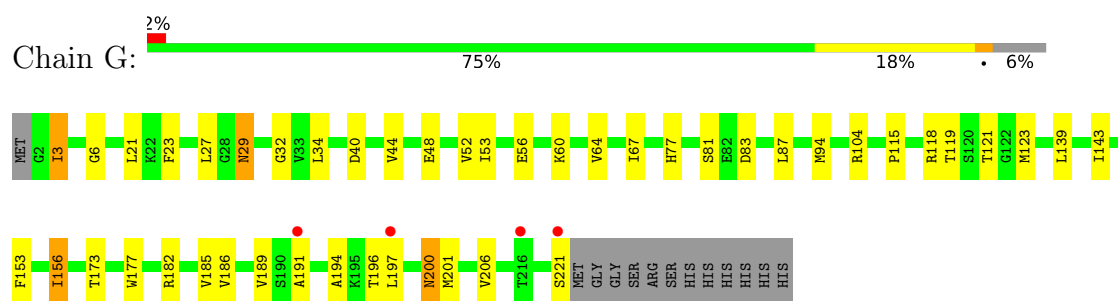
- Molecule 1: PEROXIREDOXIN 6.



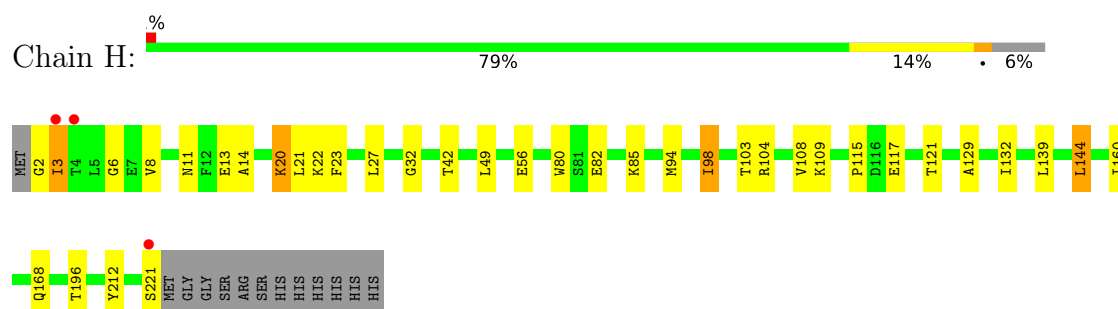
- Molecule 1: PEROXIREDOXIN 6.



- Molecule 1: PEROXIREDOXIN 6.



- Molecule 1: PEROXIREDOXIN 6.



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.11Å 111.16Å 229.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.40 15.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.8 (15.00-2.40) 97.3 (15.00-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.177 , 0.239 0.180 , 0.238	Depositor DCC
R_{free} test set	3823 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	41.2	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13959	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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: BEZ

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.16	1/1742 (0.1%)	1.14	2/2362 (0.1%)
1	B	1.12	1/1742 (0.1%)	1.07	1/2362 (0.0%)
1	C	1.14	0/1710	1.05	0/2319
1	D	1.08	2/1742 (0.1%)	1.07	4/2362 (0.2%)
1	E	1.06	3/1742 (0.2%)	1.06	3/2362 (0.1%)
1	F	1.07	2/1742 (0.1%)	1.04	1/2362 (0.0%)
1	G	1.08	3/1742 (0.2%)	1.06	2/2362 (0.1%)
1	H	1.05	2/1742 (0.1%)	1.05	1/2362 (0.0%)
All	All	1.10	14/13904 (0.1%)	1.07	14/18853 (0.1%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	205	ALA	C-O	7.02	1.32	1.24
1	G	64	VAL	CA-CB	6.83	1.62	1.54
1	F	205	ALA	CA-C	6.59	1.61	1.52
1	G	3	ILE	CA-CB	5.96	1.62	1.54
1	E	113	VAL	CA-C	5.94	1.59	1.52
1	E	113	VAL	CA-CB	5.93	1.60	1.54
1	H	160	ILE	CA-CB	5.39	1.60	1.54
1	B	156	ILE	C-O	-5.37	1.17	1.24
1	H	132	ILE	CA-CB	5.33	1.60	1.54
1	G	156	ILE	CA-CB	5.28	1.60	1.54
1	A	54	GLN	N-CA	5.24	1.52	1.46
1	D	143	ILE	CA-CB	5.09	1.62	1.55
1	D	90	VAL	CA-CB	5.02	1.59	1.53
1	E	8	VAL	C-O	-5.01	1.19	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	198	PHE	CA-C-N	6.82	126.42	119.19
1	A	198	PHE	C-N-CA	6.82	126.42	119.19
1	D	134	GLY	CA-C-N	5.85	126.25	119.47
1	D	134	GLY	C-N-CA	5.85	126.25	119.47
1	H	23	PHE	N-CA-C	5.81	117.29	111.07
1	F	23	PHE	N-CA-C	5.61	117.85	111.11
1	E	117	GLU	N-CA-C	5.45	117.12	108.67
1	G	23	PHE	N-CA-C	5.38	117.83	111.33
1	G	53	ILE	CB-CA-C	-5.38	104.88	112.14
1	B	185	VAL	N-CA-C	-5.33	101.85	108.89
1	E	216	THR	CA-C-N	-5.33	114.29	119.78
1	E	216	THR	C-N-CA	-5.33	114.29	119.78
1	D	180	GLY	N-CA-C	-5.15	107.88	114.37
1	D	23	PHE	N-CA-C	5.12	116.55	111.07

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1703	0	1723	30	0
1	B	1703	0	1723	15	0
1	C	1672	0	1690	37	0
1	D	1703	0	1723	26	0
1	E	1703	0	1723	40	0
1	F	1703	0	1723	20	0
1	G	1703	0	1723	30	0
1	H	1703	0	1723	20	0
2	A	9	0	5	0	0
2	B	9	0	5	0	0
2	C	9	0	5	1	0
2	D	9	0	5	0	0
2	E	9	0	5	1	0
2	F	9	0	5	1	0
2	G	9	0	5	0	0
2	H	9	0	5	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	46	0	0	0	0
3	B	48	0	0	0	0
3	C	33	0	0	1	0
3	D	39	0	0	1	0
3	E	31	0	0	1	0
3	F	30	0	0	1	0
3	G	33	0	0	0	0
3	H	34	0	0	1	0
All	All	13959	0	13791	189	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 7.

All (189) 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:185:VAL:HG21	1:B:201:MET:HE2	1.27	1.16
1:E:185:VAL:HG21	1:E:201:MET:HE2	1.29	1.11
1:C:34:LEU:HD12	1:C:67:ILE:HG23	1.23	1.11
1:A:185:VAL:HG21	1:A:201:MET:HE3	1.41	1.02
1:C:216:THR:HG21	1:G:121:THR:HG22	1.45	0.98
1:D:185:VAL:HG21	1:D:201:MET:CE	1.94	0.96
1:A:185:VAL:HG21	1:A:201:MET:CE	1.97	0.93
1:G:185:VAL:HG21	1:G:201:MET:HE2	1.58	0.85
1:D:34:LEU:HD13	1:D:67:ILE:HG23	1.61	0.82
1:A:171:VAL:HG11	1:A:219:PRO:HG2	1.66	0.78
1:E:57:GLY:O	1:E:61:LYS:HG2	1.85	0.77
1:C:216:THR:CG2	1:G:121:THR:HG22	2.16	0.76
1:D:185:VAL:HG21	1:D:201:MET:HE2	1.68	0.76
1:B:185:VAL:HG21	1:B:201:MET:CE	2.12	0.76
1:C:34:LEU:CD1	1:C:67:ILE:HG23	2.11	0.76
1:C:34:LEU:HD12	1:C:67:ILE:CG2	2.13	0.74
1:C:183:CYS:H	1:C:216:THR:HG22	1.53	0.73
1:D:185:VAL:HG21	1:D:201:MET:HE3	1.72	0.72
1:E:3:ILE:HG12	1:E:139:LEU:HD22	1.71	0.71
1:C:185:VAL:HG11	1:C:201:MET:CE	2.20	0.71
1:E:194:ALA:HB1	1:E:201:MET:HE1	1.73	0.70
1:C:184:MET:HE2	1:D:43:PRO:HB2	1.72	0.69
1:B:185:VAL:CG2	1:B:201:MET:HE2	2.14	0.68
1:A:40:ASP:OD2	1:A:77:HIS:HD2	1.77	0.68
1:A:60:LYS:HD3	1:H:108:VAL:HG13	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:ALA:HB1	1:A:201:MET:HE1	1.76	0.67
1:H:3:ILE:HG12	1:H:139:LEU:HD22	1.77	0.67
1:D:40:ASP:OD2	1:D:77:HIS:CD2	2.48	0.67
1:C:108:VAL:CG1	1:E:61:LYS:HD2	2.26	0.66
1:B:40:ASP:OD2	1:B:77:HIS:HD2	1.78	0.66
1:E:162:SER:O	1:E:166:THR:HG23	1.96	0.66
1:C:60:LYS:HG2	1:H:8:VAL:HB	1.79	0.65
1:A:201:MET:HE2	1:A:214:ARG:HB3	1.79	0.64
1:E:49:LEU:HD13	1:E:80:TRP:HH2	1.63	0.63
1:C:40:ASP:OD2	1:C:77:HIS:HD2	1.82	0.61
1:B:194:ALA:HB1	1:B:201:MET:HE1	1.83	0.61
1:E:67:ILE:HD12	1:E:97:PRO:HG2	1.83	0.61
1:D:143:ILE:HD12	1:D:145:TYR:CE1	2.36	0.60
1:D:67:ILE:HG12	1:D:99:ILE:HD12	1.83	0.60
1:C:182:ARG:NH2	3:C:2032:HOH:O	2.35	0.60
1:E:168:GLN:HG3	1:E:221:SER:HB2	1.84	0.60
1:G:104:ARG:NH1	1:G:123:MET:HE3	2.16	0.60
1:B:77:HIS:CE1	1:B:98:ILE:HG22	2.36	0.60
1:C:183:CYS:H	1:C:216:THR:CG2	2.13	0.60
1:E:48:GLU:O	1:E:52:VAL:HG23	2.02	0.60
1:E:40:ASP:OD2	1:E:77:HIS:HD2	1.85	0.60
1:G:40:ASP:OD2	1:G:77:HIS:HD2	1.85	0.60
1:D:109:LYS:NZ	3:D:2024:HOH:O	2.35	0.60
1:E:119:THR:HG22	1:E:125:LEU:HD21	1.83	0.59
1:D:40:ASP:OD2	1:D:77:HIS:HD2	1.84	0.59
1:C:123:MET:HE1	1:E:164:GLN:OE1	2.02	0.58
1:D:166:THR:HG22	1:D:171:VAL:O	2.03	0.58
1:E:194:ALA:HB1	1:E:201:MET:CE	2.33	0.58
1:G:29:ASN:CG	1:G:29:ASN:O	2.46	0.58
1:F:121:THR:CG2	1:F:123:MET:HG3	2.33	0.58
1:C:108:VAL:HG13	1:E:61:LYS:HD2	1.86	0.57
1:C:184:MET:HE2	1:D:43:PRO:CB	2.35	0.57
1:D:11:ASN:CG	1:D:22:LYS:NZ	2.64	0.56
1:E:81:SER:HB3	1:E:94:MET:HE3	1.88	0.56
1:E:57:GLY:O	1:E:61:LYS:CG	2.53	0.56
1:E:77:HIS:CE1	1:E:98:ILE:HG22	2.41	0.56
1:E:59:PHE:CB	1:E:66:LEU:HD11	2.36	0.56
1:G:34:LEU:HD13	1:G:67:ILE:HG23	1.87	0.56
1:A:64:VAL:HG21	1:A:160:ILE:HD13	1.88	0.55
1:A:217:PRO:O	1:A:218:GLN:C	2.49	0.55
1:D:185:VAL:CG2	1:D:201:MET:HE2	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:185:VAL:HG11	1:D:201:MET:HE1	1.88	0.55
1:C:117:GLU:CD	2:C:1218:BEZ:H4	2.32	0.54
1:F:121:THR:HG22	3:F:2015:HOH:O	2.08	0.54
1:E:198:PHE:HB2	1:E:201:MET:HE3	1.88	0.54
1:E:86:CYS:SG	1:F:207:PRO:HD3	2.48	0.53
1:F:117:GLU:CD	2:F:1222:BEZ:H4	2.33	0.53
1:E:43:PRO:HG3	1:F:212:TYR:O	2.09	0.53
1:E:59:PHE:CG	1:E:66:LEU:HD11	2.42	0.53
1:A:219:PRO:O	1:A:220:LYS:C	2.51	0.53
1:C:168:GLN:HE22	1:G:118:ARG:NE	2.07	0.52
1:C:204:LYS:HB3	1:C:213:LEU:HD23	1.92	0.52
1:E:81:SER:CB	1:E:94:MET:HE3	2.40	0.52
1:H:82:GLU:OE2	1:H:85:LYS:NZ	2.40	0.51
1:G:21:LEU:HD12	1:G:21:LEU:O	2.10	0.51
1:C:123:MET:HE1	1:E:164:GLN:CD	2.36	0.51
1:E:190:SER:HA	3:E:2026:HOH:O	2.09	0.51
1:C:216:THR:CG2	1:G:121:THR:CG2	2.87	0.51
1:C:216:THR:HG21	1:G:121:THR:CG2	2.32	0.51
1:A:174:PRO:HA	1:B:44:VAL:HG13	1.93	0.50
1:E:186:VAL:CG2	1:E:189:VAL:HG23	2.41	0.50
1:F:27:LEU:HD11	1:F:32:GLY:HA3	1.93	0.50
1:E:49:LEU:HD13	1:E:80:TRP:CH2	2.43	0.50
1:B:198:PHE:HB2	1:B:201:MET:HE3	1.94	0.49
1:H:2:GLY:N	3:H:2001:HOH:O	2.45	0.49
1:C:44:VAL:HG13	1:D:174:PRO:HA	1.93	0.49
1:H:117:GLU:HA	1:H:117:GLU:OE1	2.13	0.49
1:B:119:THR:OG1	1:B:123:MET:HB2	2.13	0.49
1:D:34:LEU:HD13	1:D:67:ILE:CG2	2.39	0.49
1:A:81:SER:O	1:A:85:LYS:HG3	2.13	0.49
1:H:11:ASN:O	1:H:22:LYS:HE2	2.12	0.48
1:F:121:THR:HG23	1:F:123:MET:HG3	1.93	0.48
1:G:21:LEU:HD12	1:G:21:LEU:C	2.37	0.48
1:A:189:VAL:HG12	1:A:194:ALA:HB2	1.95	0.48
1:C:115:PRO:HB3	1:D:6:GLY:HA3	1.95	0.48
1:C:40:ASP:OD2	1:C:77:HIS:CD2	2.64	0.47
1:G:81:SER:OG	1:G:94:MET:HE2	2.15	0.47
1:A:3:ILE:HD13	1:A:9:PHE:CD1	2.49	0.47
1:A:168:GLN:CG	1:H:121:THR:HG21	2.44	0.47
1:C:108:VAL:CG1	1:E:61:LYS:CD	2.93	0.47
1:G:153:PHE:CD1	1:G:156:ILE:HD12	2.50	0.47
1:D:11:ASN:CG	1:D:22:LYS:HZ3	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:MET:O	1:D:126:THR:HA	2.15	0.47
1:E:194:ALA:CB	1:E:201:MET:HE1	2.44	0.47
1:G:196:THR:HG22	1:G:197:LEU:HD12	1.97	0.46
1:G:115:PRO:HB3	1:H:6:GLY:HA3	1.96	0.46
1:F:49:LEU:HB3	1:F:84:VAL:HG11	1.96	0.46
1:D:27:LEU:HD11	1:D:32:GLY:HA3	1.96	0.46
1:E:57:GLY:O	1:E:61:LYS:CD	2.64	0.46
1:C:3:ILE:HG12	1:C:139:LEU:HD22	1.98	0.46
1:F:203:VAL:HG22	1:F:214:ARG:HG2	1.97	0.46
1:A:13:GLU:HG2	1:A:22:LYS:HD2	1.98	0.46
1:A:13:GLU:HA	1:A:21:LEU:O	2.17	0.45
1:A:47:THR:HG22	1:A:87:LEU:HD11	1.97	0.45
1:B:77:HIS:HE1	1:B:99:ILE:O	1.99	0.45
1:F:55:LEU:O	1:F:58:ASP:HB2	2.17	0.45
1:B:21:LEU:HD12	1:B:21:LEU:O	2.16	0.45
1:F:119:THR:OG1	1:F:121:THR:HG22	2.17	0.45
1:A:69:LEU:C	1:A:69:LEU:HD23	2.42	0.45
1:H:94:MET:CE	1:H:98:ILE:HD11	2.47	0.45
1:F:10:PRO:HB3	1:F:109:LYS:HE2	1.98	0.45
1:F:121:THR:HG21	1:F:123:MET:HG3	1.99	0.45
1:G:173:THR:HB	1:G:177:TRP:CG	2.51	0.45
1:C:14:ALA:HB1	1:C:99:ILE:HG23	1.99	0.45
1:A:48:GLU:O	1:A:52:VAL:HG23	2.18	0.44
1:B:164:GLN:O	1:B:168:GLN:CD	2.61	0.44
1:C:21:LEU:C	1:C:21:LEU:HD12	2.43	0.44
1:D:67:ILE:HG12	1:D:99:ILE:CD1	2.47	0.44
1:E:117:GLU:CD	2:E:1222:BEZ:H4	2.42	0.44
1:G:194:ALA:CB	1:G:201:MET:HE1	2.48	0.44
1:A:153:PHE:HA	1:A:156:ILE:HD12	1.99	0.44
1:G:27:LEU:HD11	1:G:32:GLY:HA3	1.99	0.44
1:C:34:LEU:HD11	1:C:99:ILE:HD12	2.00	0.44
1:G:3:ILE:HG12	1:G:139:LEU:HD22	1.99	0.44
1:G:119:THR:OG1	1:G:123:MET:HB2	2.17	0.44
1:C:162:SER:O	1:C:166:THR:HG23	2.18	0.43
1:G:6:GLY:HA3	1:H:115:PRO:HB3	1.99	0.43
1:G:191:ALA:O	1:G:194:ALA:HB3	2.18	0.43
1:E:6:GLY:HA3	1:F:115:PRO:HB3	2.00	0.43
1:F:68:ALA:HB3	1:F:98:ILE:HG12	2.00	0.43
1:C:168:GLN:NE2	1:G:118:ARG:NE	2.66	0.43
1:H:27:LEU:HD11	1:H:32:GLY:HA3	1.99	0.43
1:F:121:THR:HG23	1:F:123:MET:CG	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:THR:HA	1:A:87:LEU:CD1	2.48	0.43
1:A:185:VAL:HG21	1:A:201:MET:HE2	1.93	0.43
1:D:26:TRP:CH2	1:D:67:ILE:HB	2.54	0.42
1:E:59:PHE:HB2	1:E:66:LEU:HD11	2.00	0.42
1:G:29:ASN:O	1:G:29:ASN:OD1	2.36	0.42
1:H:49:LEU:HD13	1:H:80:TRP:HH2	1.84	0.42
1:B:194:ALA:HB1	1:B:201:MET:CE	2.47	0.42
1:D:153:PHE:HA	1:D:156:ILE:HD12	2.02	0.42
1:H:42:THR:HB	2:H:1222:BEZ:C	2.50	0.42
1:A:57:GLY:O	1:A:61:LYS:HG3	2.20	0.42
1:D:52:VAL:CG1	1:D:66:LEU:HD13	2.50	0.42
1:E:14:ALA:HB1	1:E:99:ILE:HG23	2.02	0.42
1:H:21:LEU:C	1:H:21:LEU:HD12	2.44	0.42
1:A:220:LYS:HA	1:A:220:LYS:HD2	1.95	0.42
1:G:200:ASN:HD22	1:G:200:ASN:HA	1.74	0.42
1:A:168:GLN:HG3	1:H:121:THR:HG21	2.00	0.42
1:D:185:VAL:CG2	1:D:201:MET:CE	2.82	0.42
1:F:114:ASP:CG	1:F:115:PRO:HD2	2.45	0.41
1:F:220:LYS:O	1:F:220:LYS:NZ	2.33	0.41
1:C:206:VAL:O	1:C:207:PRO:C	2.63	0.41
1:E:216:THR:HG23	1:E:217:PRO:O	2.20	0.41
1:C:108:VAL:HG11	1:E:61:LYS:HG3	2.02	0.41
1:B:48:GLU:O	1:B:52:VAL:HG23	2.21	0.41
1:C:174:PRO:HG2	1:C:182:ARG:NH2	2.36	0.41
1:G:44:VAL:O	1:G:48:GLU:HG3	2.20	0.41
1:H:103:THR:O	1:H:104:ARG:HB2	2.20	0.41
1:A:93:ASP:O	1:A:94:MET:C	2.62	0.41
1:C:69:LEU:HD23	1:C:69:LEU:C	2.46	0.41
1:A:6:GLY:HA3	1:B:115:PRO:HB3	2.03	0.41
1:A:182:ARG:HB3	1:A:217:PRO:HA	2.03	0.41
1:C:108:VAL:HG11	1:E:61:LYS:CD	2.51	0.40
1:G:83:ASP:OD2	1:H:212:TYR:OH	2.36	0.40
1:A:26:TRP:CH2	1:A:65:LYS:HB3	2.56	0.40
1:E:197:LEU:HD23	1:E:198:PHE:CE2	2.56	0.40
1:F:49:LEU:O	1:F:96:TYR:OH	2.38	0.40
1:G:48:GLU:O	1:G:52:VAL:HG23	2.22	0.40
1:G:186:VAL:O	1:G:189:VAL:HG12	2.21	0.40
1:H:14:ALA:O	1:H:20:LYS:HA	2.21	0.40
1:E:36:SER:HA	1:E:69:LEU:O	2.21	0.40
1:F:40:ASP:OD2	1:F:77:HIS:CD2	2.74	0.40
1:H:129:ALA:HB2	1:H:144:LEU:HD12	2.03	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	218/233 (94%)	209 (96%)	9 (4%)	0	100	100
1	B	218/233 (94%)	211 (97%)	7 (3%)	0	100	100
1	C	214/233 (92%)	205 (96%)	9 (4%)	0	100	100
1	D	218/233 (94%)	213 (98%)	5 (2%)	0	100	100
1	E	218/233 (94%)	211 (97%)	6 (3%)	1 (0%)	24	37
1	F	218/233 (94%)	212 (97%)	6 (3%)	0	100	100
1	G	218/233 (94%)	206 (94%)	12 (6%)	0	100	100
1	H	218/233 (94%)	210 (96%)	8 (4%)	0	100	100
All	All	1740/1864 (93%)	1677 (96%)	62 (4%)	1 (0%)	48	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	43	PRO

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	189/200 (94%)	184 (97%)	5 (3%)	40	63

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	189/200 (94%)	185 (98%)	4 (2%)	47	69
1	C	185/200 (92%)	176 (95%)	9 (5%)	22	39
1	D	189/200 (94%)	181 (96%)	8 (4%)	26	45
1	E	189/200 (94%)	175 (93%)	14 (7%)	13	22
1	F	189/200 (94%)	180 (95%)	9 (5%)	23	40
1	G	189/200 (94%)	180 (95%)	9 (5%)	23	40
1	H	189/200 (94%)	179 (95%)	10 (5%)	20	36
All	All	1508/1600 (94%)	1440 (96%)	68 (4%)	24	42

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	GLU
1	A	21	LEU
1	A	87	LEU
1	A	94	MET
1	A	144	LEU
1	B	22	LYS
1	B	34	LEU
1	B	87	LEU
1	B	195	LYS
1	C	87	LEU
1	C	98	ILE
1	C	123	MET
1	C	143	ILE
1	C	182	ARG
1	C	197	LEU
1	C	204	LYS
1	C	206	VAL
1	C	216	THR
1	D	21	LEU
1	D	34	LEU
1	D	39	ARG
1	D	82	GLU
1	D	144	LEU
1	D	196	THR
1	D	197	LEU
1	D	201	MET
1	E	3	ILE
1	E	20	LYS

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Mol	Chain	Res	Type
1	E	30	SER
1	E	49	LEU
1	E	56	GLU
1	E	61	LYS
1	E	67	ILE
1	E	72	ASP
1	E	86	CYS
1	E	87	LEU
1	E	143	ILE
1	E	144	LEU
1	E	197	LEU
1	E	216	THR
1	F	29	ASN
1	F	87	LEU
1	F	108	VAL
1	F	120	SER
1	F	123	MET
1	F	143	ILE
1	F	168	GLN
1	F	202	GLU
1	F	220	LYS
1	G	29	ASN
1	G	56	GLU
1	G	60	LYS
1	G	87	LEU
1	G	143	ILE
1	G	182	ARG
1	G	200	ASN
1	G	206	VAL
1	G	221	SER
1	H	3	ILE
1	H	13	GLU
1	H	20	LYS
1	H	56	GLU
1	H	98	ILE
1	H	109	LYS
1	H	144	LEU
1	H	168	GLN
1	H	196	THR
1	H	221	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	GLN
1	A	178	GLN
1	B	29	ASN
1	B	77	HIS
1	B	178	GLN
1	C	77	HIS
1	C	168	GLN
1	C	178	GLN
1	D	77	HIS
1	D	178	GLN
1	E	77	HIS
1	E	178	GLN
1	F	168	GLN
1	F	178	GLN
1	G	77	HIS
1	G	178	GLN
1	G	200	ASN
1	H	73	ASN
1	H	77	HIS
1	H	178	GLN

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 ⓘ

8 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	BEZ	C	1218	-	9,9,9	0.99	0	11,11,11	1.17	1 (9%)
2	BEZ	B	1222	-	9,9,9	0.92	0	11,11,11	1.58	4 (36%)
2	BEZ	H	1222	-	9,9,9	0.71	0	11,11,11	1.55	3 (27%)
2	BEZ	F	1222	-	9,9,9	0.96	0	11,11,11	1.88	4 (36%)
2	BEZ	G	1222	-	9,9,9	0.98	1 (11%)	11,11,11	1.68	4 (36%)
2	BEZ	A	1222	-	9,9,9	0.92	0	11,11,11	1.54	2 (18%)
2	BEZ	D	1222	-	9,9,9	1.05	0	11,11,11	1.58	2 (18%)
2	BEZ	E	1222	-	9,9,9	0.85	0	11,11,11	1.86	3 (27%)

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	BEZ	C	1218	-	-	0/4/4/4	0/1/1/1
2	BEZ	B	1222	-	-	0/4/4/4	0/1/1/1
2	BEZ	H	1222	-	-	0/4/4/4	0/1/1/1
2	BEZ	F	1222	-	-	4/4/4/4	0/1/1/1
2	BEZ	G	1222	-	-	1/4/4/4	0/1/1/1
2	BEZ	A	1222	-	-	1/4/4/4	0/1/1/1
2	BEZ	D	1222	-	-	4/4/4/4	0/1/1/1
2	BEZ	E	1222	-	-	0/4/4/4	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	1222	BEZ	O2-C	-2.77	1.22	1.30

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1222	BEZ	O2-C-C1	3.51	123.84	114.84
2	E	1222	BEZ	C6-C1-C2	3.18	122.61	118.57
2	A	1222	BEZ	C3-C2-C1	-3.09	117.32	120.36
2	F	1222	BEZ	C3-C2-C1	-3.08	117.34	120.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1222	BEZ	O2-C-C1	3.03	122.62	114.84
2	A	1222	BEZ	C6-C1-C2	2.94	122.30	118.57
2	F	1222	BEZ	C6-C1-C2	2.83	122.17	118.57
2	B	1222	BEZ	C4-C3-C2	-2.81	116.77	120.24
2	F	1222	BEZ	O2-C-C1	2.80	122.03	114.84
2	G	1222	BEZ	O2-C-C1	2.75	121.90	114.84
2	H	1222	BEZ	C5-C6-C1	-2.61	117.80	120.36
2	G	1222	BEZ	C4-C3-C2	-2.52	117.13	120.24
2	G	1222	BEZ	O2-C-O1	-2.51	117.96	123.35
2	C	1218	BEZ	O2-C-C1	2.46	121.16	114.84
2	D	1222	BEZ	O2-C-O1	-2.41	118.18	123.35
2	H	1222	BEZ	O2-C-C1	2.40	120.99	114.84
2	E	1222	BEZ	O1-C-C1	-2.38	115.21	121.46
2	H	1222	BEZ	C6-C1-C2	2.24	121.41	118.57
2	B	1222	BEZ	C3-C2-C1	2.18	122.51	120.36
2	B	1222	BEZ	C5-C4-C3	2.17	122.84	119.87
2	F	1222	BEZ	C4-C5-C6	-2.16	117.57	120.24
2	B	1222	BEZ	O2-C-O1	-2.09	118.85	123.35
2	G	1222	BEZ	C5-C6-C1	-2.05	118.34	120.36

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	1222	BEZ	O1-C-C1-C2
2	F	1222	BEZ	O1-C-C1-C6
2	F	1222	BEZ	O2-C-C1-C2
2	F	1222	BEZ	O2-C-C1-C6
2	D	1222	BEZ	O2-C-C1-C2
2	D	1222	BEZ	O2-C-C1-C6
2	D	1222	BEZ	O1-C-C1-C2
2	D	1222	BEZ	O1-C-C1-C6
2	G	1222	BEZ	O1-C-C1-C2
2	A	1222	BEZ	O1-C-C1-C2

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1218	BEZ	1	0
2	H	1222	BEZ	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1222	BEZ	1	0
2	E	1222	BEZ	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	220/233 (94%)	-0.20	2 (0%) 81 78	34, 40, 46, 62	0
1	B	220/233 (94%)	-0.21	0 100 100	33, 40, 47, 58	0
1	C	216/233 (92%)	-0.13	1 (0%) 87 85	34, 41, 45, 50	0
1	D	220/233 (94%)	-0.16	1 (0%) 87 85	34, 40, 47, 58	0
1	E	220/233 (94%)	0.53	4 (1%) 67 63	34, 41, 47, 53	0
1	F	220/233 (94%)	0.51	5 (2%) 61 57	35, 41, 47, 51	0
1	G	220/233 (94%)	0.02	4 (1%) 67 63	35, 41, 47, 49	0
1	H	220/233 (94%)	0.27	3 (1%) 73 69	35, 40, 46, 52	0
All	All	1756/1864 (94%)	0.08	20 (1%) 78 74	33, 41, 47, 62	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	221	SER	3.8
1	C	182	ARG	2.8
1	E	75	ALA	2.7
1	E	148	THR	2.7
1	G	197	LEU	2.6
1	A	221	SER	2.4
1	D	221	SER	2.4
1	G	221	SER	2.4
1	F	208	SER	2.3
1	E	114	ASP	2.3
1	F	29	ASN	2.3
1	F	120	SER	2.2
1	H	3	ILE	2.2
1	E	88	SER	2.2
1	G	191	ALA	2.1
1	F	221	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	216	THR	2.1
1	H	4	THR	2.0
1	F	83	ASP	2.0
1	A	219	PRO	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](#)

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. 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	B-factors($\text{\AA}^2$)	Q<0.9
2	BEZ	F	1222	9/9	0.91	0.13	47,48,49,49	0
2	BEZ	H	1222	9/9	0.91	0.14	39,39,40,40	0
2	BEZ	D	1222	9/9	0.92	0.14	31,33,35,36	0
2	BEZ	E	1222	9/9	0.92	0.14	42,42,43,43	0
2	BEZ	C	1218	9/9	0.93	0.14	36,37,38,38	0
2	BEZ	A	1222	9/9	0.95	0.13	25,28,29,29	0
2	BEZ	B	1222	9/9	0.96	0.14	29,30,32,32	0
2	BEZ	G	1222	9/9	0.97	0.12	30,33,33,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.