



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 2, 2016 – 01:57 AM EDT

PDB ID : 5B6N
Title : Crystal structures of human peroxiredoxin 6 in sulfinic acid state
Authors : Kim, K.H.; Lee, W.T.; Kim, E.E.
Deposited on : 2016-05-30
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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
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) : rb-20027939

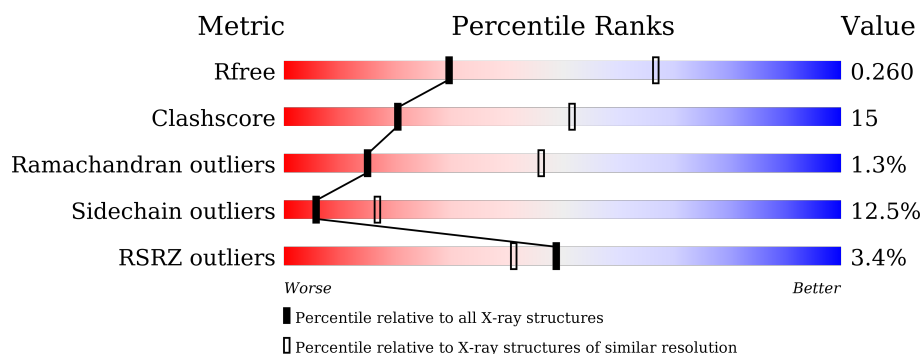
1 Overall quality at a glance ⓘ

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	224	<div> <div>68%</div> <div>27%</div> <div>..</div> </div>
1	B	224	<div> <div>3%</div> <div>67%</div> <div>28%</div> <div>..</div> </div>
1	C	224	<div> <div>4%</div> <div>64%</div> <div>26%</div> <div>7%</div> <div>.</div> </div>
1	D	224	<div> <div>10%</div> <div>49%</div> <div>39%</div> <div>7%</div> <div>5%</div> </div>
1	E	224	<div> <div>4%</div> <div>56%</div> <div>37%</div> <div>6%</div> <div>.</div> </div>
1	F	224	<div> <div>60%</div> <div>32%</div> <div>8%</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
1	CSD	E	47	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10576 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	221	Total	C	N	O	S	0	0	0
			1749	1127	293	324	5			
1	B	221	Total	C	N	O	S	0	0	0
			1749	1127	293	324	5			
1	C	217	Total	C	N	O	S	0	0	0
			1723	1111	288	319	5			
1	D	213	Total	C	N	O	S	0	0	0
			1690	1092	285	309	4			
1	E	221	Total	C	N	O	S	0	0	0
			1749	1127	293	324	5			
1	F	223	Total	C	N	O	S	0	0	0
			1760	1134	295	326	5			

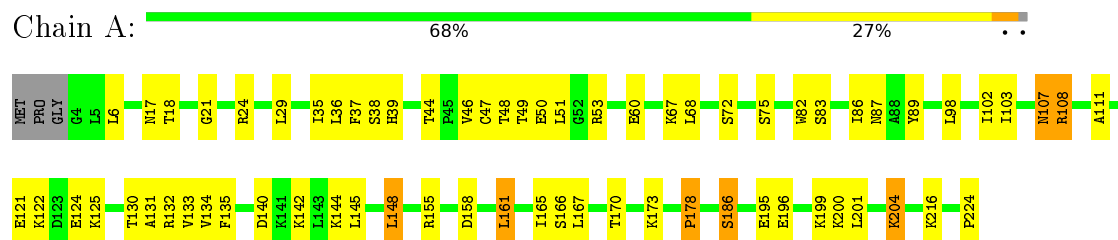
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	28	Total	O	0	0
			28	28		
2	B	22	Total	O	0	0
			22	22		
2	C	29	Total	O	0	0
			29	29		
2	D	21	Total	O	0	0
			21	21		
2	E	27	Total	O	0	0
			27	27		
2	F	29	Total	O	0	0
			29	29		

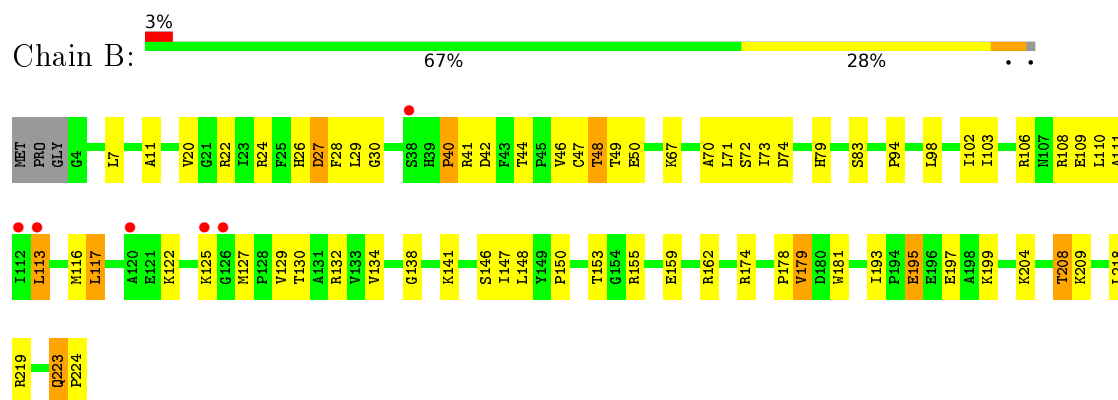
3 Residue-property plots [i](#)

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.

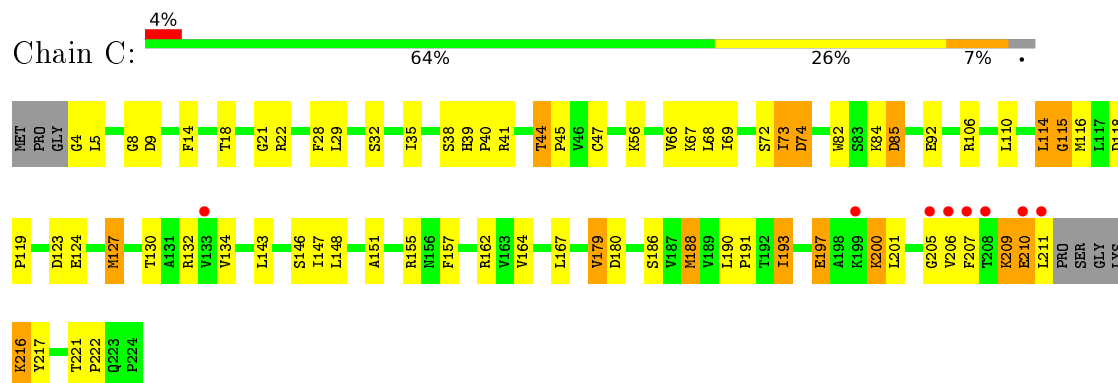
• Molecule 1: Peroxiredoxin-6



• Molecule 1: Peroxiredoxin-6

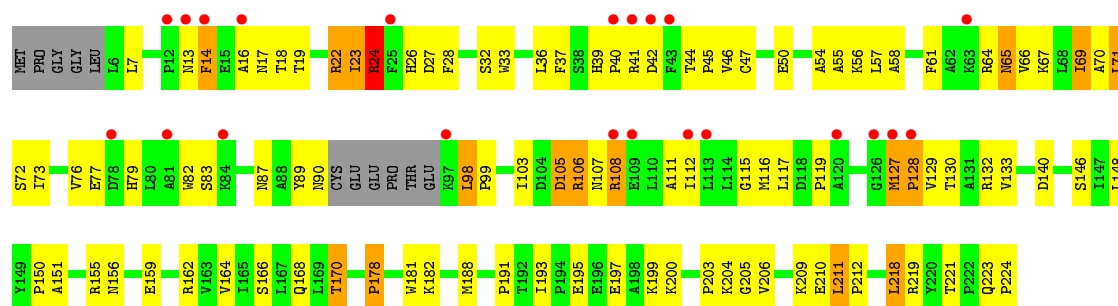


• Molecule 1: Peroxiredoxin-6

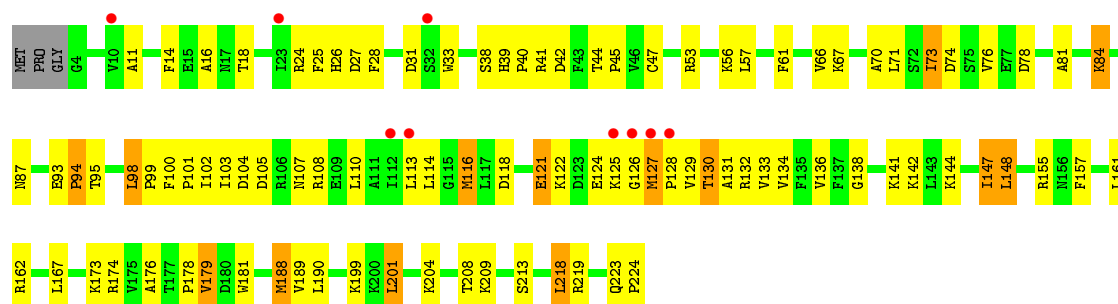


• Molecule 1: Peroxiredoxin-6

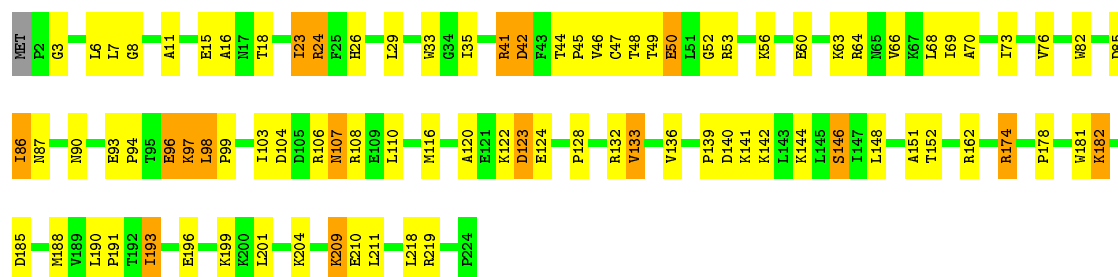




• Molecule 1: Peroxiredoxin-6



• 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	94.42Å 106.54Å 167.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.99 – 2.90 30.18 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.8 (29.99-2.90) 97.9 (30.18-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.90Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.248 , 0.304 0.248 , 0.260	Depositor DCC
R_{free} test set	1886 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.145	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 14.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	10576	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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: CSD

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	0.64	0/1783	0.55	0/2420
1	B	0.62	0/1783	0.54	0/2420
1	C	0.59	0/1755	0.58	0/2381
1	D	0.51	0/1722	0.59	1/2335 (0.0%)
1	E	0.55	0/1783	0.57	0/2420
1	F	0.59	0/1795	0.56	0/2436
All	All	0.58	0/10621	0.57	1/14412 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	24	ARG	N-CA-C	-5.51	96.12	111.00

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	1749	0	1778	43	1
1	B	1749	0	1778	48	0
1	C	1723	0	1749	49	0
1	D	1690	0	1726	76	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1749	0	1778	59	0
1	F	1760	0	1788	57	1
2	A	28	0	0	0	0
2	B	22	0	0	0	0
2	C	29	0	0	0	0
2	D	21	0	0	0	0
2	E	27	0	0	0	0
2	F	29	0	0	0	0
All	All	10576	0	10597	307	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (307) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:PHE:O	1:C:67:LYS:NZ	2.10	0.85
1:D:108:ARG:HG2	1:D:128:PRO:HB2	1.59	0.83
1:C:41:ARG:O	1:C:44:THR:OG1	1.99	0.81
1:A:204:LYS:HE3	1:A:204:LYS:H	1.49	0.78
1:B:44:THR:O	1:B:48:THR:OG1	2.02	0.78
1:E:39:HIS:NE2	1:E:42:ASP:OD2	2.16	0.77
1:D:205:GLY:O	1:D:221:THR:OG1	2.03	0.74
1:B:208:THR:HB	1:B:219:ARG:HG2	1.68	0.74
1:A:83:SER:O	1:A:87:ASN:ND2	2.20	0.73
1:B:174:ARG:NH1	1:B:197:GLU:OE1	2.22	0.73
1:C:197:GLU:O	1:C:200:LYS:NZ	2.21	0.73
1:F:96:GLU:O	1:F:98:LEU:N	2.20	0.73
1:C:186:SER:HA	1:C:222:PRO:HA	1.70	0.72
1:F:193:ILE:O	1:F:219:ARG:NH2	2.18	0.72
1:F:35:ILE:HG13	1:F:136:VAL:HG12	1.71	0.71
1:A:6:LEU:HD13	1:B:117:LEU:HD21	1.72	0.71
1:A:50:GLU:HG2	1:B:179:VAL:HG13	1.72	0.71
1:E:162:ARG:HD3	1:E:181:TRP:O	1.91	0.70
1:D:32:SER:HB3	1:D:65:ASN:HD21	1.56	0.70
1:A:166:SER:O	1:A:170:THR:OG1	2.08	0.70
1:A:38:SER:OG	1:A:131:ALA:O	2.09	0.70
1:A:37:PHE:HB2	1:A:155:ARG:HH12	1.56	0.69
1:D:106:ARG:NH1	1:D:107:ASN:OD1	2.25	0.69
1:D:111:ALA:O	1:D:115:GLY:N	2.24	0.68
1:E:11:ALA:O	1:E:26:HIS:NE2	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:CSD:OD1	1:B:132:ARG:NH1	2.27	0.68
1:A:44:THR:O	1:A:48:THR:OG1	2.13	0.67
1:F:33:TRP:HB2	1:F:66:VAL:HG22	1.75	0.67
1:F:52:GLY:O	1:F:90:ASN:ND2	2.27	0.66
1:C:167:LEU:HD21	1:D:150:PRO:HG3	1.78	0.66
1:D:159:GLU:OE1	1:D:162:ARG:NE	2.29	0.66
1:B:11:ALA:O	1:B:26:HIS:NE2	2.22	0.66
1:E:130:THR:OG1	1:E:131:ALA:N	2.27	0.66
1:D:18:THR:OG1	1:D:19:THR:N	2.25	0.65
1:E:47:CSD:OD2	1:E:132:ARG:NH1	2.28	0.65
1:F:162:ARG:NH1	1:F:181:TRP:O	2.26	0.65
1:C:18:THR:OG1	1:C:21:GLY:O	2.13	0.65
1:E:208:THR:HG23	1:E:219:ARG:HG2	1.77	0.65
1:C:72:SER:OG	1:C:73:ILE:N	2.22	0.65
1:D:33:TRP:HB2	1:D:65:ASN:O	1.97	0.64
1:E:38:SER:OG	1:E:131:ALA:O	2.16	0.64
1:D:72:SER:HB3	1:D:79:HIS:HE1	1.60	0.64
1:B:71:LEU:HD12	1:B:103:ILE:HD12	1.80	0.64
1:C:56:LYS:NZ	1:C:92:GLU:OE2	2.28	0.64
1:E:129:VAL:HG12	1:E:130:THR:H	1.63	0.62
1:D:61:PHE:HD1	1:D:66:VAL:HG21	1.64	0.62
1:F:116:MET:HB2	1:F:133:VAL:HG11	1.81	0.62
1:A:18:THR:OG1	1:A:21:GLY:O	2.18	0.61
1:E:122:LYS:HB3	1:E:126:GLY:HA2	1.81	0.61
1:E:39:HIS:HB2	1:E:47:CSD:HB2	1.83	0.61
1:E:47:CSD:HD2	1:E:132:ARG:HH12	1.48	0.61
1:E:31:ASP:HA	1:E:141:LYS:HE3	1.82	0.61
1:F:139:PRO:O	1:F:141:LYS:NZ	2.34	0.61
1:D:47:CSD:N	1:D:47:CSD:OD2	2.34	0.60
1:A:39:HIS:NE2	1:A:72:SER:HB2	2.16	0.60
1:C:35:ILE:HD11	1:C:164:VAL:HG21	1.83	0.60
1:D:54:ALA:HA	1:D:57:LEU:HD12	1.83	0.60
1:A:47:CSD:N	1:A:47:CSD:OD2	2.33	0.60
1:A:29:LEU:HA	1:A:67:LYS:HD2	1.83	0.60
1:A:145:LEU:HD12	1:B:150:PRO:HD3	1.82	0.60
1:D:156:ASN:OD1	1:D:159:GLU:N	2.35	0.60
1:F:48:THR:HG23	1:F:86:ILE:HG22	1.83	0.60
1:E:44:THR:HG23	1:F:191:PRO:HG3	1.84	0.60
1:A:133:VAL:HG22	1:A:148:LEU:HG	1.83	0.59
1:B:46:VAL:O	1:B:49:THR:OG1	2.15	0.59
1:C:188:MET:SD	1:C:188:MET:N	2.75	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:LEU:HD23	1:D:99:PRO:HD2	1.83	0.59
1:F:18:THR:HG22	1:F:103:ILE:HG23	1.83	0.59
1:F:120:ALA:O	1:F:122:LYS:NZ	2.24	0.59
1:A:196:GLU:HA	1:A:199:LYS:HE2	1.85	0.59
1:F:151:ALA:O	1:F:152:THR:OG1	2.18	0.58
1:B:150:PRO:HD2	1:B:153:THR:HB	1.85	0.58
1:F:11:ALA:O	1:F:26:HIS:NE2	2.30	0.58
1:B:30:GLY:O	1:B:141:LYS:NZ	2.36	0.58
1:E:40:PRO:HA	1:E:130:THR:HG23	1.85	0.58
1:A:167:LEU:HD21	1:B:150:PRO:HG2	1.84	0.58
1:B:42:ASP:OD2	1:B:79:HIS:ND1	2.36	0.58
1:A:36:LEU:HD23	1:A:135:PHE:HD2	1.67	0.57
1:A:148:LEU:HD23	1:B:7:LEU:HG	1.87	0.57
1:C:134:VAL:HB	1:C:147:ILE:HG12	1.86	0.57
1:D:69:ILE:HD11	1:D:103:ILE:HG13	1.85	0.57
1:D:129:VAL:HG22	1:D:130:THR:H	1.69	0.57
1:B:108:ARG:CZ	1:B:127:MET:HB3	2.34	0.57
1:F:35:ILE:HB	1:F:68:LEU:HD23	1.85	0.56
1:D:41:ARG:HG2	1:D:42:ASP:H	1.70	0.56
1:D:79:HIS:O	1:D:83:SER:N	2.31	0.56
1:E:61:PHE:HB3	1:E:66:VAL:HB	1.87	0.56
1:D:18:THR:HG1	1:D:19:THR:H	1.51	0.56
1:D:24:ARG:HG3	1:D:26:HIS:HB2	1.87	0.56
1:F:41:ARG:NH2	1:F:123:ASP:HB2	2.21	0.55
1:F:53:ARG:HH11	1:F:56:LYS:HE3	1.70	0.55
1:C:8:GLY:HA3	1:D:119:PRO:HG2	1.88	0.55
1:D:64:ARG:O	1:D:66:VAL:N	2.39	0.55
1:C:179:VAL:HG13	1:D:50:GLU:HG2	1.89	0.55
1:D:58:ALA:HA	1:D:61:PHE:HB2	1.89	0.55
1:F:122:LYS:HA	1:F:128:PRO:HA	1.89	0.55
1:F:136:VAL:HG23	1:F:144:LYS:HB2	1.89	0.55
1:A:178:PRO:HA	1:B:46:VAL:HG13	1.88	0.54
1:D:193:ILE:O	1:D:219:ARG:NH2	2.40	0.54
1:E:70:ALA:HB3	1:E:102:ILE:HG23	1.89	0.54
1:E:71:LEU:HA	1:E:103:ILE:HG22	1.89	0.54
1:F:8:GLY:O	1:F:142:LYS:HE3	2.07	0.54
1:A:46:VAL:HG13	1:B:178:PRO:HA	1.88	0.54
1:F:70:ALA:O	1:F:103:ILE:HG13	2.07	0.54
1:E:174:ARG:HD3	1:E:201:LEU:HD11	1.88	0.54
1:A:46:VAL:O	1:A:49:THR:OG1	2.22	0.54
1:D:22:ARG:HA	1:D:22:ARG:CZ	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:ASP:N	1:A:140:ASP:OD1	2.36	0.54
1:C:47:CSD:OD2	1:C:132:ARG:NH1	2.41	0.53
1:D:42:ASP:HA	1:D:82:TRP:CZ3	2.43	0.53
1:A:53:ARG:NH2	1:A:158:ASP:OD1	2.39	0.53
1:D:18:THR:HG22	1:D:23:ILE:HG21	1.91	0.53
1:A:37:PHE:HB2	1:A:155:ARG:NH1	2.23	0.53
1:E:121:GLU:O	1:E:129:VAL:HG23	2.09	0.53
1:F:53:ARG:HA	1:F:56:LYS:HE2	1.90	0.53
1:A:108:ARG:HG2	1:A:111:ALA:HB3	1.91	0.53
1:C:191:PRO:HG3	1:D:45:PRO:HD2	1.90	0.53
1:E:41:ARG:HH12	1:E:124:GLU:HG3	1.74	0.53
1:B:111:ALA:HA	1:B:116:MET:HE2	1.91	0.52
1:B:79:HIS:O	1:B:83:SER:OG	2.26	0.52
1:D:16:ALA:H	1:D:23:ILE:HD11	1.74	0.52
1:C:216:LYS:HG3	1:C:217:TYR:HD2	1.74	0.52
1:C:29:LEU:O	1:C:32:SER:OG	2.27	0.52
1:E:189:VAL:HB	1:E:219:ARG:HB2	1.90	0.52
1:E:73:ILE:HG13	1:E:108:ARG:HG2	1.91	0.52
1:A:35:ILE:HB	1:A:68:LEU:HD23	1.90	0.52
1:D:197:GLU:HA	1:D:200:LYS:HD2	1.91	0.52
1:E:133:VAL:HG22	1:E:148:LEU:HD23	1.90	0.52
1:A:140:ASP:O	1:A:142:LYS:HG3	2.10	0.51
1:C:134:VAL:HG22	1:C:155:ARG:HH11	1.76	0.51
1:F:60:GLU:OE1	1:F:64:ARG:NH1	2.43	0.51
1:E:81:ALA:HA	1:E:84:LYS:HD3	1.92	0.51
1:C:5:LEU:HD11	1:C:114:LEU:HD23	1.93	0.51
1:D:50:GLU:OE2	1:D:155:ARG:HG2	2.11	0.51
1:D:133:VAL:N	1:D:155:ARG:HH12	2.07	0.51
1:E:138:GLY:HA3	1:E:144:LYS:NZ	2.26	0.51
1:E:178:PRO:HA	1:F:46:VAL:HG13	1.91	0.51
1:F:45:PRO:O	1:F:49:THR:HG23	2.10	0.51
1:D:162:ARG:HD2	1:D:181:TRP:O	2.10	0.51
1:D:111:ALA:HB1	1:D:117:LEU:HB2	1.91	0.50
1:C:35:ILE:HB	1:C:68:LEU:HD23	1.94	0.50
1:D:36:LEU:HD11	1:D:71:LEU:HD13	1.94	0.50
1:F:86:ILE:HD11	1:F:98:LEU:HD11	1.94	0.50
1:A:38:SER:O	1:A:155:ARG:NH2	2.38	0.50
1:F:29:LEU:HB2	1:F:141:LYS:HE3	1.94	0.50
1:D:178:PRO:HG3	1:D:188:MET:SD	2.52	0.50
1:E:84:LYS:HB3	1:E:95:THR:HA	1.92	0.50
1:A:173:LYS:HG3	1:A:224:PRO:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:193:ILE:O	1:B:219:ARG:NH2	2.37	0.49
1:B:40:PRO:HD2	1:B:47:CSD:SG	2.52	0.49
1:D:64:ARG:HB3	1:D:168:GLN:HE22	1.76	0.49
1:E:148:LEU:HD13	1:F:7:LEU:HG	1.95	0.49
1:E:121:GLU:O	1:E:128:PRO:HA	2.12	0.49
1:D:18:THR:HB	1:D:103:ILE:HA	1.95	0.49
1:B:108:ARG:NH2	1:B:127:MET:HB3	2.27	0.49
1:D:73:ILE:HD12	1:D:106:ARG:HA	1.95	0.49
1:D:14:PHE:HD1	1:D:23:ILE:HD12	1.77	0.49
1:E:116:MET:HB3	1:E:133:VAL:HG21	1.95	0.49
1:C:197:GLU:HG2	1:C:200:LYS:HZ2	1.78	0.48
1:D:105:ASP:C	1:D:107:ASN:H	2.15	0.48
1:D:166:SER:O	1:D:170:THR:OG1	2.17	0.48
1:D:223:GLN:HA	1:D:224:PRO:HD3	1.71	0.48
1:F:93:GLU:H	1:F:97:LYS:HZ1	1.60	0.48
1:B:44:THR:OG1	1:B:47:CSD:OD2	2.14	0.48
1:D:197:GLU:O	1:D:200:LYS:HB2	2.14	0.47
1:F:174:ARG:HG3	1:F:201:LEU:HD21	1.97	0.47
1:B:108:ARG:NH2	1:B:127:MET:O	2.48	0.47
1:B:72:SER:OG	1:B:73:ILE:N	2.47	0.47
1:D:111:ALA:HA	1:D:116:MET:HB2	1.95	0.47
1:C:66:VAL:HG21	1:C:164:VAL:HG11	1.95	0.47
1:F:82:TRP:O	1:F:86:ILE:HG23	2.14	0.47
1:E:47:CSD:HD2	1:E:132:ARG:NH1	2.09	0.47
1:A:107:ASN:HD22	1:A:107:ASN:H	1.62	0.47
1:A:144:LYS:O	1:B:150:PRO:HG3	2.14	0.47
1:B:159:GLU:OE2	1:B:162:ARG:NH2	2.31	0.47
1:B:209:LYS:HB3	1:B:218:LEU:HD23	1.97	0.47
1:F:47:CSD:SG	1:F:132:ARG:NH1	2.83	0.47
1:E:148:LEU:N	1:F:146:SER:O	2.34	0.47
1:D:57:LEU:O	1:D:61:PHE:N	2.47	0.47
1:E:213:SER:OG	1:F:85:ASP:OD2	2.20	0.47
1:B:71:LEU:HG	1:B:103:ILE:HB	1.97	0.46
1:B:73:ILE:HD12	1:B:108:ARG:HG3	1.96	0.46
1:C:82:TRP:O	1:C:85:ASP:N	2.42	0.46
1:C:38:SER:O	1:C:155:ARG:NH2	2.46	0.46
1:B:110:LEU:HA	1:B:113:LEU:HB2	1.96	0.46
1:C:148:LEU:HD12	1:D:146:SER:OG	2.15	0.46
1:C:14:PHE:CG	1:C:110:LEU:HD21	2.51	0.46
1:E:110:LEU:HA	1:E:113:LEU:HB2	1.97	0.46
1:E:84:LYS:HG2	1:E:84:LYS:H	1.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:108:ARG:H	1:D:108:ARG:HD2	1.81	0.46
1:E:179:VAL:CG1	1:F:50:GLU:HG3	2.46	0.46
1:E:93:GLU:HA	1:E:94:PRO:HA	1.75	0.46
1:C:114:LEU:O	1:C:116:MET:N	2.49	0.46
1:B:28:PHE:O	1:B:67:LYS:NZ	2.38	0.45
1:C:74:ASP:OD2	1:C:74:ASP:N	2.49	0.45
1:E:136:VAL:HG23	1:E:144:LYS:HB2	1.98	0.45
1:D:14:PHE:CD1	1:D:23:ILE:HD12	2.51	0.45
1:A:18:THR:HG22	1:A:103:ILE:HA	1.98	0.45
1:B:26:HIS:O	1:B:30:GLY:N	2.46	0.45
1:F:193:ILE:HG12	1:F:193:ILE:H	1.41	0.45
1:D:203:PRO:HG2	1:D:204:LYS:HE2	1.99	0.45
1:F:15:GLU:HA	1:F:23:ILE:O	2.17	0.45
1:C:197:GLU:HG2	1:C:200:LYS:NZ	2.32	0.44
1:F:96:GLU:HG3	1:F:96:GLU:H	1.61	0.44
1:B:134:VAL:O	1:B:146:SER:HA	2.17	0.44
1:C:45:PRO:HD3	1:D:191:PRO:HB3	1.99	0.44
1:C:132:ARG:NH2	1:C:151:ALA:O	2.44	0.44
1:F:15:GLU:HB3	1:F:24:ARG:HD2	2.00	0.44
1:F:93:GLU:HA	1:F:94:PRO:HD3	1.89	0.44
1:E:33:TRP:CE3	1:E:138:GLY:HA2	2.53	0.44
1:F:69:ILE:HG12	1:F:103:ILE:HD11	2.00	0.44
1:F:107:ASN:N	1:F:107:ASN:OD1	2.51	0.44
1:A:161:LEU:O	1:A:165:ILE:HG13	2.17	0.44
1:D:46:VAL:O	1:D:132:ARG:NH2	2.51	0.44
1:E:45:PRO:HB3	1:F:218:LEU:HA	1.99	0.44
1:C:134:VAL:CG2	1:C:155:ARG:HH11	2.31	0.44
1:E:223:GLN:HA	1:E:224:PRO:HD3	1.76	0.44
1:B:41:ARG:HD2	1:B:74:ASP:OD2	2.18	0.44
1:C:39:HIS:NE2	1:C:72:SER:HB2	2.33	0.44
1:B:70:ALA:HB3	1:B:102:ILE:HG23	2.00	0.43
1:D:56:LYS:NZ	1:D:90:ASN:HA	2.32	0.43
1:E:24:ARG:HG2	1:E:27:ASP:HB2	2.00	0.43
1:D:108:ARG:NH1	1:D:127:MET:HB3	2.33	0.43
1:D:178:PRO:HD2	1:D:181:TRP:CD1	2.53	0.43
1:D:211:LEU:HD21	1:D:218:LEU:HD12	1.99	0.43
1:A:49:THR:HB	1:B:179:VAL:HG22	2.01	0.43
1:C:5:LEU:HB3	1:C:143:LEU:HD22	1.98	0.43
1:D:55:ALA:C	1:D:57:LEU:H	2.21	0.43
1:C:190:LEU:O	1:C:193:ILE:HG23	2.18	0.43
1:C:84:LYS:O	1:D:212:PRO:HG2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:76:VAL:HG22	1:F:104:ASP:HB2	2.00	0.43
1:B:110:LEU:HD23	1:B:110:LEU:H	1.83	0.43
1:F:209:LYS:HE3	1:F:210:GLU:N	2.34	0.43
1:C:72:SER:OG	1:C:74:ASP:OD2	2.35	0.43
1:F:41:ARG:NH2	1:F:124:GLU:HG2	2.33	0.43
1:C:179:VAL:HB	1:C:180:ASP:OD1	2.19	0.43
1:C:47:CSD:OD1	1:C:47:CSD:N	2.52	0.43
1:B:22:ARG:NH2	1:C:66:VAL:O	2.52	0.43
1:F:23:ILE:H	1:F:23:ILE:HG13	1.53	0.43
1:C:4:GLY:HA3	1:C:115:GLY:H	1.83	0.43
1:D:209:LYS:O	1:D:211:LEU:HD22	2.18	0.43
1:F:44:THR:OG1	1:F:47:CSD:OD2	2.05	0.42
1:A:98:LEU:HD12	1:A:102:ILE:HD11	2.01	0.42
1:C:209:LYS:HE3	1:C:209:LYS:HA	2.01	0.42
1:E:18:THR:HG22	1:E:103:ILE:HG13	2.01	0.42
1:E:98:LEU:HA	1:E:99:PRO:HD3	1.85	0.42
1:F:42:ASP:O	1:F:82:TRP:CD2	2.72	0.42
1:D:72:SER:OG	1:D:73:ILE:N	2.52	0.42
1:E:134:VAL:HG23	1:E:147:ILE:HD11	2.02	0.42
1:D:132:ARG:HH11	1:D:151:ALA:HA	1.84	0.42
1:D:204:LYS:HD3	1:D:204:LYS:HA	1.71	0.42
1:B:195:GLU:O	1:B:199:LYS:HE2	2.20	0.42
1:F:182:LYS:NZ	1:F:185:ASP:OD2	2.38	0.42
1:C:110:LEU:O	1:C:114:LEU:HD12	2.20	0.42
1:E:199:LYS:HB3	1:E:199:LYS:HE2	1.71	0.42
1:B:223:GLN:HA	1:B:224:PRO:HD3	1.86	0.42
1:E:100:PHE:HA	1:E:101:PRO:HD3	1.87	0.42
1:B:162:ARG:HD2	1:B:181:TRP:O	2.19	0.42
1:E:44:THR:HB	1:E:47:CSD:SG	2.60	0.42
1:F:56:LYS:HE2	1:F:56:LYS:HB2	1.77	0.42
1:A:51:LEU:HD22	1:A:82:TRP:CZ3	2.54	0.41
1:D:41:ARG:HG2	1:D:42:ASP:N	2.34	0.41
1:E:28:PHE:CZ	1:E:67:LYS:HB3	2.55	0.41
1:D:195:GLU:O	1:D:199:LYS:HG2	2.20	0.41
1:E:76:VAL:HG23	1:E:104:ASP:HB2	2.02	0.41
1:A:134:VAL:CG2	1:A:155:ARG:HH11	2.32	0.41
1:A:134:VAL:HG22	1:A:155:ARG:HH11	1.84	0.41
1:E:127:MET:SD	1:E:128:PRO:HD2	2.60	0.41
1:F:16:ALA:HB2	1:F:110:LEU:HD11	2.01	0.41
1:F:98:LEU:HA	1:F:99:PRO:HD2	1.78	0.41
1:E:132:ARG:HB3	1:E:155:ARG:NE	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:190:LEU:HD12	1:E:190:LEU:H	1.84	0.41
1:A:124:GLU:HG2	1:A:125:LYS:HG3	2.03	0.41
1:D:50:GLU:OE2	1:D:155:ARG:N	2.53	0.41
1:D:76:VAL:HA	1:D:79:HIS:HD2	1.85	0.41
1:A:132:ARG:HB3	1:A:155:ARG:NE	2.35	0.41
1:B:134:VAL:HB	1:B:147:ILE:HG12	2.03	0.41
1:C:123:ASP:N	1:C:127:MET:O	2.47	0.41
1:C:200:LYS:HZ3	1:C:201:LEU:HB2	1.85	0.41
1:A:195:GLU:CD	1:A:195:GLU:H	2.23	0.41
1:A:47:CSD:OD2	1:A:132:ARG:NH1	2.44	0.41
1:F:140:ASP:OD2	1:F:142:LYS:HB2	2.21	0.41
1:F:73:ILE:HG13	1:F:73:ILE:H	1.62	0.41
1:D:39:HIS:NE2	1:D:72:SER:HB2	2.36	0.41
1:B:138:GLY:O	1:B:141:LYS:N	2.45	0.41
1:C:210:GLU:O	1:C:211:LEU:HD23	2.20	0.41
1:D:33:TRP:HB3	1:D:66:VAL:HG12	2.03	0.41
1:E:144:LYS:HB3	1:E:167:LEU:HD13	2.03	0.41
1:B:125:LYS:HA	1:B:125:LYS:HD2	1.98	0.41
1:A:86:ILE:O	1:A:89:TYR:HB3	2.20	0.40
1:D:211:LEU:HD21	1:D:218:LEU:HB2	2.01	0.40
1:D:72:SER:HB3	1:D:79:HIS:CE1	2.48	0.40
1:D:115:GLY:C	1:D:117:LEU:H	2.24	0.40
1:D:42:ASP:HA	1:D:82:TRP:CE3	2.57	0.40
1:E:16:ALA:HB3	1:E:103:ILE:HD11	2.03	0.40
1:E:209:LYS:HB3	1:E:218:LEU:HB3	2.03	0.40
1:F:106:ARG:O	1:F:108:ARG:NH1	2.51	0.40
1:F:87:ASN:HD22	1:F:87:ASN:HA	1.74	0.40
1:B:27:ASP:OD2	1:B:27:ASP:N	2.52	0.40
1:C:119:PRO:HG2	1:D:7:LEU:O	2.20	0.40
1:D:37:PHE:O	1:D:70:ALA:HA	2.21	0.40
1:E:118:ASP:HB3	1:E:121:GLU:HB2	2.03	0.40
1:B:49:THR:OG1	1:B:50:GLU:N	2.54	0.40
1:C:118:ASP:HA	1:C:119:PRO:HD3	1.85	0.40
1:C:200:LYS:NZ	1:C:201:LEU:HB2	2.37	0.40
1:E:176:ALA:O	1:E:188:MET:HB2	2.21	0.40
1:E:57:LEU:HD23	1:E:57:LEU:HA	1.97	0.40

All (1) 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:186:SER:OG	1:F:204:LYS:O[1_655]	2.13	0.07

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	218/224 (97%)	204 (94%)	13 (6%)	1 (0%)	34	71
1	B	218/224 (97%)	199 (91%)	17 (8%)	2 (1%)	21	57
1	C	212/224 (95%)	186 (88%)	23 (11%)	3 (1%)	14	44
1	D	208/224 (93%)	171 (82%)	32 (15%)	5 (2%)	7	29
1	E	218/224 (97%)	185 (85%)	30 (14%)	3 (1%)	14	44
1	F	220/224 (98%)	193 (88%)	24 (11%)	3 (1%)	14	44
All	All	1294/1344 (96%)	1138 (88%)	139 (11%)	17 (1%)	15	46

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	97	LYS
1	C	205	GLY
1	D	40	PRO
1	D	65	ASN
1	C	115	GLY
1	D	128	PRO
1	E	125	LYS
1	F	178	PRO
1	E	25	PHE
1	E	94	PRO
1	D	140	ASP
1	F	3	GLY
1	A	178	PRO
1	B	94	PRO
1	D	178	PRO

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Mol	Chain	Res	Type
1	C	40	PRO
1	B	40	PRO

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	191/193 (99%)	175 (92%)	16 (8%)	14	37
1	B	191/193 (99%)	171 (90%)	20 (10%)	8	25
1	C	188/193 (97%)	162 (86%)	26 (14%)	4	13
1	D	184/193 (95%)	155 (84%)	29 (16%)	3	9
1	E	191/193 (99%)	164 (86%)	27 (14%)	4	12
1	F	192/193 (100%)	168 (88%)	24 (12%)	6	17
All	All	1137/1158 (98%)	995 (88%)	142 (12%)	6	17

All (142) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	24	ARG
1	A	60	GLU
1	A	75	SER
1	A	107	ASN
1	A	108	ARG
1	A	121	GLU
1	A	122	LYS
1	A	130	THR
1	A	148	LEU
1	A	161	LEU
1	A	186	SER
1	A	200	LYS
1	A	201	LEU
1	A	204	LYS
1	A	216	LYS

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Mol	Chain	Res	Type
1	B	20	VAL
1	B	24	ARG
1	B	27	ASP
1	B	29	LEU
1	B	48	THR
1	B	98	LEU
1	B	106	ARG
1	B	109	GLU
1	B	113	LEU
1	B	117	LEU
1	B	122	LYS
1	B	129	VAL
1	B	130	THR
1	B	148	LEU
1	B	155	ARG
1	B	179	VAL
1	B	195	GLU
1	B	204	LYS
1	B	208	THR
1	B	223	GLN
1	C	9	ASP
1	C	22	ARG
1	C	44	THR
1	C	69	ILE
1	C	73	ILE
1	C	74	ASP
1	C	85	ASP
1	C	106	ARG
1	C	114	LEU
1	C	124	GLU
1	C	127	MET
1	C	130	THR
1	C	146	SER
1	C	157	PHE
1	C	162	ARG
1	C	179	VAL
1	C	188	MET
1	C	193	ILE
1	C	197	GLU
1	C	200	LYS
1	C	206	VAL
1	C	207	PHE

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Mol	Chain	Res	Type
1	C	209	LYS
1	C	210	GLU
1	C	216	LYS
1	C	221	THR
1	D	13	ASN
1	D	14	PHE
1	D	17	ASN
1	D	22	ARG
1	D	23	ILE
1	D	24	ARG
1	D	27	ASP
1	D	28	PHE
1	D	44	THR
1	D	67	LYS
1	D	69	ILE
1	D	71	LEU
1	D	77	GLU
1	D	87	ASN
1	D	89	TYR
1	D	98	LEU
1	D	105	ASP
1	D	106	ARG
1	D	108	ARG
1	D	112	ILE
1	D	127	MET
1	D	148	LEU
1	D	164	VAL
1	D	170	THR
1	D	182	LYS
1	D	206	VAL
1	D	210	GLU
1	D	211	LEU
1	D	218	LEU
1	E	14	PHE
1	E	53	ARG
1	E	56	LYS
1	E	73	ILE
1	E	74	ASP
1	E	78	ASP
1	E	84	LYS
1	E	87	ASN
1	E	98	LEU

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Mol	Chain	Res	Type
1	E	105	ASP
1	E	107	ASN
1	E	114	LEU
1	E	116	MET
1	E	121	GLU
1	E	127	MET
1	E	130	THR
1	E	142	LYS
1	E	147	ILE
1	E	148	LEU
1	E	157	PHE
1	E	161	LEU
1	E	173	LYS
1	E	179	VAL
1	E	188	MET
1	E	201	LEU
1	E	204	LYS
1	E	218	LEU
1	F	6	LEU
1	F	23	ILE
1	F	24	ARG
1	F	41	ARG
1	F	42	ASP
1	F	50	GLU
1	F	63	LYS
1	F	86	ILE
1	F	96	GLU
1	F	98	LEU
1	F	107	ASN
1	F	123	ASP
1	F	133	VAL
1	F	146	SER
1	F	148	LEU
1	F	174	ARG
1	F	182	LYS
1	F	188	MET
1	F	190	LEU
1	F	193	ILE
1	F	196	GLU
1	F	199	LYS
1	F	209	LYS
1	F	211	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	ASN
1	B	223	GLN
1	D	65	ASN
1	D	79	HIS
1	F	65	ASN
1	F	87	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CSD	A	47	1	2,7,8	0.69	0	2,8,10	0.69	0
1	CSD	B	47	1	2,7,8	0.73	0	2,8,10	0.70	0
1	CSD	C	47	1	2,7,8	0.78	0	2,8,10	1.21	0
1	CSD	D	47	1	2,7,8	0.75	0	2,8,10	0.68	0
1	CSD	E	47	1	2,7,8	0.74	0	2,8,10	0.60	0
1	CSD	F	47	1	2,7,8	0.67	0	2,8,10	1.41	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
1	CSD	A	47	1	-	0/2/6/8	0/0/0/0
1	CSD	B	47	1	-	0/2/6/8	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSD	C	47	1	-	0/2/6/8	0/0/0/0
1	CSD	D	47	1	-	1/2/6/8	0/0/0/0
1	CSD	E	47	1	-	1/2/6/8	0/0/0/0
1	CSD	F	47	1	-	1/2/6/8	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	D	47	CSD	CA-CB-SG-OD1
1	E	47	CSD	CA-CB-SG-OD1
1	F	47	CSD	CA-CB-SG-OD1

There are no ring outliers.

6 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	47	CSD	2	0
1	B	47	CSD	3	0
1	C	47	CSD	2	0
1	D	47	CSD	1	0
1	E	47	CSD	5	0
1	F	47	CSD	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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 ⓘ

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	220/224 (98%)	-0.31	0 100 100	0, 9, 29, 44	0
1	B	220/224 (98%)	-0.10	6 (2%) 58 52	0, 11, 33, 59	0
1	C	216/224 (96%)	-0.13	8 (3%) 45 38	0, 13, 46, 59	0
1	D	212/224 (94%)	0.46	22 (10%) 8 5	2, 33, 64, 71	0
1	E	220/224 (98%)	0.18	9 (4%) 41 34	1, 24, 54, 76	0
1	F	222/224 (99%)	-0.28	0 100 100	0, 15, 35, 64	0
All	All	1310/1344 (97%)	-0.03	45 (3%) 49 41	0, 16, 52, 76	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	120	ALA	7.1
1	D	112	ILE	5.8
1	D	41	ARG	5.3
1	E	126	GLY	4.6
1	D	14	PHE	4.4
1	D	12	PRO	4.1
1	D	108	ARG	4.1
1	D	78	ASP	3.9
1	D	13	ASN	3.9
1	E	128	PRO	3.9
1	D	126	GLY	3.6
1	C	208	THR	3.6
1	B	112	ILE	3.5
1	E	112	ILE	3.5
1	E	127	MET	3.5
1	C	205	GLY	3.3
1	D	127	MET	3.3
1	C	210	GLU	3.2
1	E	10	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	211	LEU	3.2
1	D	128	PRO	3.2
1	C	207	PHE	3.0
1	D	42	ASP	2.9
1	C	206	VAL	2.9
1	D	109	GLU	2.9
1	B	113	LEU	2.8
1	D	81	ALA	2.8
1	E	125	LYS	2.7
1	D	113	LEU	2.5
1	D	16	ALA	2.4
1	E	113	LEU	2.4
1	B	38	SER	2.4
1	D	43	PHE	2.4
1	D	84	LYS	2.4
1	C	133	VAL	2.3
1	C	199	LYS	2.3
1	D	63	LYS	2.3
1	B	126	GLY	2.2
1	D	120	ALA	2.2
1	E	23	ILE	2.2
1	D	97	LYS	2.2
1	E	32	SER	2.1
1	D	25	PHE	2.1
1	D	40	PRO	2.0
1	B	125	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	CSD	F	47	8/9	0.97	0.22	-	11,11,11,11	0
1	CSD	B	47	8/9	0.94	0.32	-	15,15,15,15	0
1	CSD	C	47	8/9	0.96	0.20	-	8,8,8,8	0
1	CSD	E	47	8/9	0.95	0.30	-	16,16,16,16	0
1	CSD	A	47	8/9	0.93	0.18	-	5,5,5,5	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	CSD	D	47	8/9	0.92	0.25	-	43,43,43,43	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.