



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 06:10 PM UTC

PDB ID : 6HF1 / pdb_00006hf1
Title : Mutant oxidoreductase fragment of mouse QSOX1 in complex with an anti-body Fab
Authors : Grossman-Haham, I.; Fass, D.
Deposited on : 2018-08-21
Resolution : 1.94 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
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

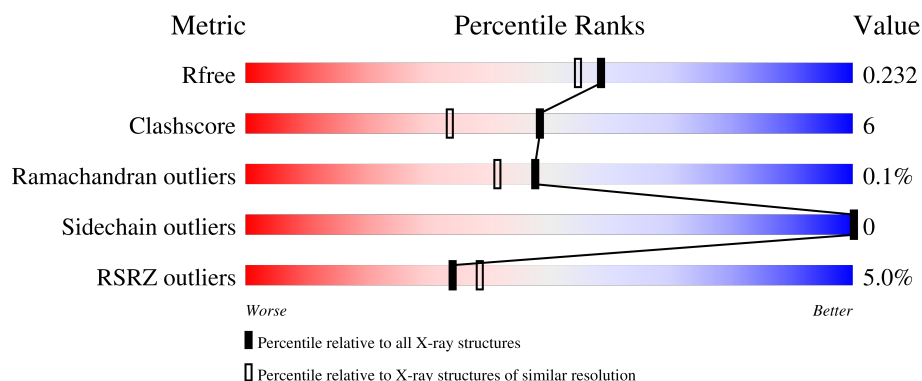
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	238	10% 86% 14%
1	D	238	5% 86% 14%
2	C	215	6% 88% 9% .
3	B	212	6% 86% 13%
3	E	212	92% 7%

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Mol	Chain	Length	Quality of chain
4	F	216	2% 88% 10%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10954 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 Sulfhydryl oxidase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	238	Total	C	N	O	S	0	2	0
			1854	1179	318	350	7			
1	D	238	Total	C	N	O	S	0	0	0
			1847	1173	318	349	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	75	ALA	HIS	engineered mutation	UNP Q8BND5
A	122	THR	PRO	engineered mutation	UNP Q8BND5
D	75	ALA	HIS	engineered mutation	UNP Q8BND5
D	122	THR	PRO	engineered mutation	UNP Q8BND5

- Molecule 2 is a protein called Fab 316 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	209	Total	C	N	O	S	0	2	0
			1603	1024	261	313	5			

- Molecule 3 is a protein called Fab 316 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	211	Total	C	N	O	S	0	1	0
			1616	1015	267	327	7			
3	E	212	Total	C	N	O	S	0	2	0
			1627	1022	269	329	7			

- Molecule 4 is a protein called Fab 316 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	210	Total	C	N	O	S	0	1	0
			1603	1024	261	313	5			

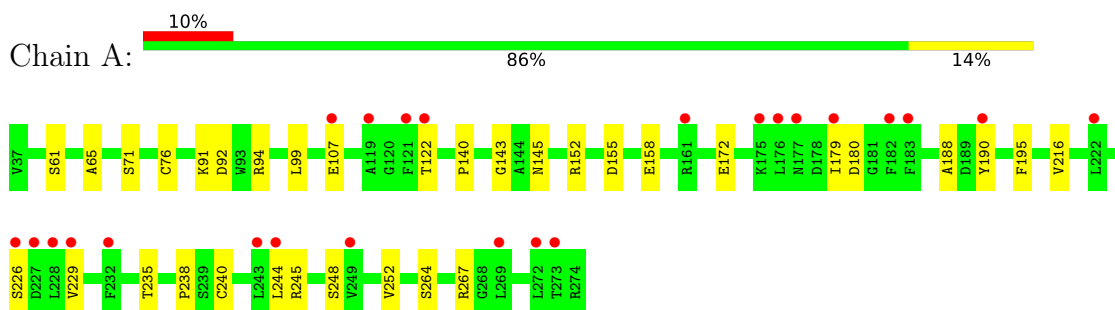
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	126	Total 126	O 126	0	0
5	C	133	Total 133	O 133	0	0
5	B	163	Total 163	O 163	0	0
5	D	101	Total 101	O 101	0	0
5	F	155	Total 155	O 155	0	0
5	E	126	Total 126	O 126	0	0

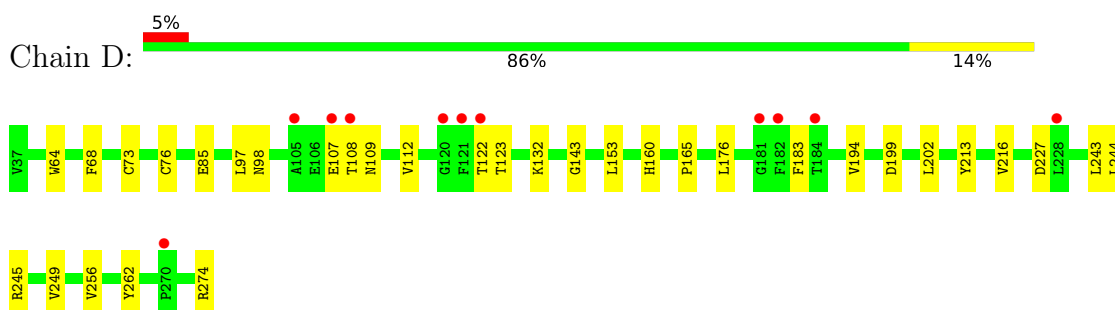
3 Residue-property plots

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.

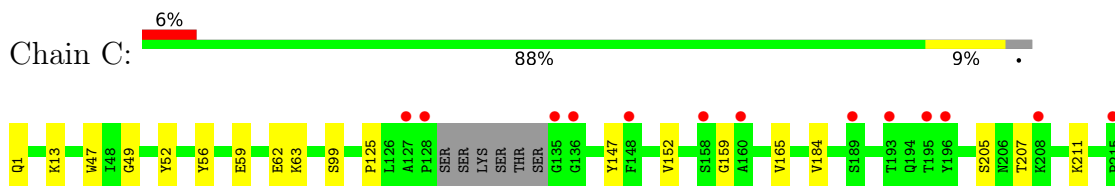
• Molecule 1: Sulfhydryl oxidase 1



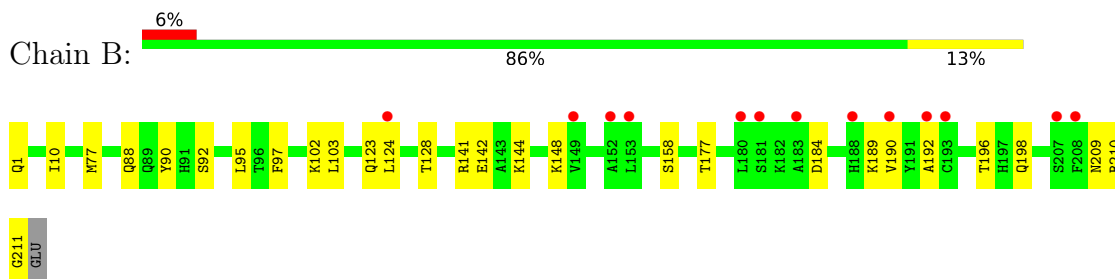
• Molecule 1: Sulfhydryl oxidase 1



• Molecule 2: Fab 316 heavy chain



• Molecule 3: Fab 316 light chain

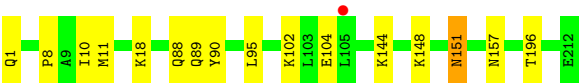


● Molecule 3: Fab 316 light chain

Chain E:

92%

7%



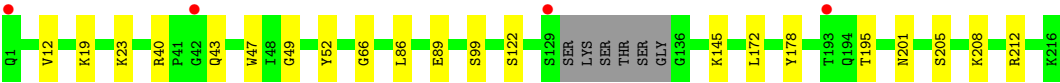
● Molecule 4: Fab 316 heavy chain

Chain F:

2%

88%

10%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.54Å 112.72Å 193.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.89 – 1.94 48.89 – 1.94	Depositor EDS
% Data completeness (in resolution range)	97.6 (48.89-1.94) 98.0 (48.89-1.94)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 1.94Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.194 , 0.231 0.195 , 0.232	Depositor DCC
R_{free} test set	5215 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.282	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10954	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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](#)

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	0.31	0/1904	0.52	0/2598
1	D	0.31	0/1891	0.49	0/2580
2	C	0.32	0/1654	0.53	0/2258
3	B	0.30	0/1659	0.51	0/2255
3	E	0.31	0/1673	0.56	0/2274
4	F	0.31	0/1651	0.52	0/2255
All	All	0.31	0/10432	0.52	0/14220

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	C	0	1
4	F	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	99	SER	Peptide
4	F	99	SER	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1854	0	1815	29	1
1	D	1847	0	1801	25	0
2	C	1603	0	1559	15	2
3	B	1616	0	1564	24	0
3	E	1627	0	1574	10	0
4	F	1603	0	1552	17	0
5	A	126	0	0	10	1
5	B	163	0	0	4	0
5	C	133	0	0	4	0
5	D	101	0	0	3	0
5	E	126	0	0	3	0
5	F	155	0	0	6	0
All	All	10954	0	9865	110	2

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 6.

All (110) 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)
3:B:10:ILE:HD13	3:B:102:LYS:HB3	1.34	1.06
3:B:144:LYS:NZ	5:B:301:HOH:O	1.92	1.02
1:A:61:SER:O	5:A:302:HOH:O	1.83	0.95
3:B:1:GLN:OE1	3:B:1:GLN:N	2.09	0.86
3:B:10:ILE:HD12	3:B:102:LYS:O	1.80	0.82
1:A:158:GLU:OE2	5:A:303:HOH:O	1.99	0.81
3:B:10:ILE:HD13	3:B:102:LYS:CB	2.13	0.75
3:B:10:ILE:CD1	3:B:102:LYS:HB3	2.15	0.75
1:D:107:GLU:N	1:D:107:GLU:OE1	2.20	0.75
1:A:179:ILE:HB	5:A:311:HOH:O	1.87	0.74
4:F:43:GLN:OE1	5:F:301:HOH:O	2.06	0.73
1:A:172:GLU:OE2	5:A:304:HOH:O	2.08	0.71
1:A:71:SER:O	5:A:305:HOH:O	2.09	0.69
1:D:243:LEU:HG	1:D:249:VAL:HG22	1.74	0.69
1:A:180:ASP:N	5:A:311:HOH:O	2.25	0.68
1:D:85:GLU:OE2	5:D:301:HOH:O	2.11	0.68
4:F:23:LYS:HD2	5:F:336:HOH:O	1.93	0.68
2:C:159:GLY:O	5:C:301:HOH:O	2.11	0.67
1:A:235:THR:O	5:A:306:HOH:O	2.14	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:PRO:O	5:A:307:HOH:O	2.14	0.65
3:B:1:GLN:H1	3:B:1:GLN:CD	2.02	0.65
3:E:148:LYS:HB3	3:E:151:ASN:HA	1.80	0.63
1:D:213:TYR:OH	1:D:274:ARG:NH1	2.27	0.62
2:C:62:GLU:OE2	5:C:302:HOH:O	2.16	0.62
1:D:183:PHE:O	1:D:245:ARG:NH2	2.33	0.61
1:A:122:THR:OG1	5:A:308:HOH:O	2.16	0.60
1:D:68:PHE:O	1:D:122:THR:HA	2.02	0.60
4:F:40[A]:ARG:NE	5:F:301:HOH:O	2.29	0.60
3:E:1:GLN:OE1	3:E:1:GLN:N	2.30	0.60
1:A:91:LYS:HD2	1:A:94:ARG:HD2	1.85	0.59
1:A:91:LYS:HE3	1:A:94:ARG:HH11	1.68	0.57
1:D:216:VAL:O	5:D:302:HOH:O	2.18	0.57
1:A:65:ALA:HB3	1:A:99:LEU:HD22	1.87	0.57
1:D:107:GLU:HG2	1:D:108:THR:N	2.18	0.56
1:A:152[B]:ARG:NH1	1:A:155:ASP:OD2	2.38	0.56
2:C:13:LYS:HE3	1:D:256:VAL:HG13	1.87	0.56
2:C:125:PRO:HD3	2:C:211:LYS:HE2	1.88	0.55
1:A:92:ASP:OD2	1:A:267:ARG:NH2	2.40	0.55
1:A:76:CYS:SG	1:A:122:THR:HB	2.47	0.54
1:A:190:TYR:HB2	1:A:216:VAL:HG22	1.88	0.54
3:B:148:LYS:HB2	3:B:192:ALA:HB3	1.88	0.54
4:F:66:GLY:N	5:F:302:HOH:O	2.09	0.53
3:B:141:ARG:NH1	5:B:311:HOH:O	2.41	0.53
3:E:10:ILE:HD11	3:E:104:GLU:OE1	2.09	0.53
1:A:244:LEU:HD23	1:A:248:SER:OG	2.08	0.53
1:A:188:ALA:O	1:A:245:ARG:NH1	2.43	0.52
3:E:157:ASN:HB2	5:E:361:HOH:O	2.08	0.52
1:A:226:SER:HA	1:A:229:VAL:HB	1.92	0.51
3:B:198:GLN:NE2	5:B:302:HOH:O	2.19	0.51
2:C:63:LYS:NZ	5:C:304:HOH:O	2.32	0.51
1:A:195:PHE:O	1:A:238:PRO:HA	2.12	0.49
1:D:97:LEU:HD21	1:D:153:LEU:HD13	1.95	0.49
1:D:176:LEU:HD22	1:D:227:ASP:HB3	1.95	0.49
3:E:8:PRO:HG2	3:E:11:MET:HB3	1.95	0.48
3:E:18:LYS:NZ	5:E:301:HOH:O	2.25	0.48
1:D:107:GLU:HG2	1:D:108:THR:H	1.78	0.48
1:D:109:ASN:O	1:D:112:VAL:HG12	2.13	0.47
1:A:143:GLY:HA3	2:C:52:TYR:CD2	2.49	0.47
1:A:145:ASN:HB2	2:C:59:GLU:OE2	2.14	0.47
3:E:10:ILE:HD13	3:E:102:LYS:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:202:LEU:HD21	1:D:262:TYR:CE2	2.50	0.47
2:C:1:GLN:N	5:C:307:HOH:O	2.41	0.46
3:E:90:TYR:HA	3:E:95:LEU:HD22	1.97	0.46
4:F:12:VAL:HG21	4:F:86:LEU:HD13	1.97	0.46
4:F:145:LYS:HE2	5:E:304:HOH:O	2.15	0.46
4:F:195:THR:HG23	4:F:212:ARG:HD2	1.98	0.46
1:D:160:HIS:CD2	1:D:165:PRO:HD3	2.51	0.46
3:B:184:ASP:OD1	3:B:184:ASP:N	2.49	0.46
1:A:143:GLY:HA3	2:C:52:TYR:CG	2.51	0.45
1:A:264[B]:SER:HB2	4:F:205:SER:HB2	1.98	0.45
2:C:147:TYR:CE2	2:C:152:VAL:HG13	2.52	0.45
4:F:122:SER:OG	5:F:303:HOH:O	2.21	0.45
3:B:210:ARG:HA	3:B:211:GLY:HA2	1.63	0.45
1:A:264[A]:SER:HB3	4:F:205:SER:HB2	1.99	0.44
3:B:10:ILE:HD13	3:B:102:LYS:CG	2.46	0.44
2:C:13:LYS:HE3	1:D:256:VAL:CG1	2.48	0.44
1:D:122:THR:O	1:D:123:THR:OG1	2.28	0.44
3:B:92:SER:HB3	5:B:400:HOH:O	2.17	0.44
3:B:144:LYS:HB3	3:B:196:THR:HB	1.99	0.44
1:A:240:CYS:HB3	1:A:252:VAL:HB	2.00	0.43
1:A:152[B]:ARG:HA	1:A:152[B]:ARG:HD3	1.69	0.43
3:B:190:VAL:HG22	3:B:209:ASN:OD1	2.19	0.43
1:D:143:GLY:HA3	4:F:52:TYR:CG	2.53	0.43
3:B:124:LEU:HD23	3:B:124:LEU:HA	1.84	0.43
4:F:172:LEU:HD13	4:F:178:TYR:CZ	2.54	0.43
1:A:140:PRO:O	1:A:152[A]:ARG:NH1	2.52	0.42
3:B:88:GLN:HG3	3:B:97:PHE:CE2	2.54	0.42
3:E:88:GLN:NE2	3:E:89:GLN:O	2.53	0.42
1:D:73:CYS:HB3	1:D:76:CYS:SG	2.59	0.42
3:B:90:TYR:HA	3:B:95:LEU:HD22	2.01	0.42
1:A:248:SER:HA	5:A:313:HOH:O	2.20	0.42
3:B:77:MET:SD	3:B:103:LEU:HD21	2.59	0.42
2:C:165:VAL:HG22	2:C:184:VAL:HB	2.02	0.42
3:B:189:LYS:HG2	3:B:190:VAL:HG23	2.02	0.42
1:D:143:GLY:HA3	4:F:52:TYR:CD2	2.55	0.41
1:D:194:VAL:HG11	1:D:202:LEU:HD23	2.02	0.41
3:B:158:SER:HA	3:B:177:THR:O	2.20	0.41
1:D:64:TRP:CZ3	1:D:98:ASN:HB3	2.55	0.41
3:B:123:GLN:HG2	3:B:128:THR:O	2.20	0.41
3:B:142:GLU:H	3:B:142:GLU:CD	2.28	0.41
2:C:63:LYS:HE2	4:F:89:GLU:CD	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:244:LEU:HA	1:D:244:LEU:HD12	1.70	0.41
4:F:47:TRP:CZ2	4:F:49:GLY:HA2	2.56	0.41
2:C:205:SER:O	2:C:207:THR:HG23	2.21	0.41
3:E:144:LYS:HB3	3:E:196:THR:HB	2.02	0.41
2:C:47:TRP:CZ2	2:C:49:GLY:HA2	2.56	0.41
4:F:19:LYS:NZ	5:F:315:HOH:O	2.53	0.40
1:D:132:LYS:NZ	5:D:306:HOH:O	2.38	0.40
1:D:199:ASP:OD1	1:D:199:ASP:N	2.54	0.40
4:F:201:ASN:ND2	4:F:208:LYS:HG2	2.36	0.40

All (2) 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)
2:C:56:TYR:OH	5:A:301:HOH:O[4_555]	1.93	0.27
1:A:107:GLU:OE1	2:C:56:TYR:OH[4_455]	2.14	0.06

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	238/238 (100%)	231 (97%)	7 (3%)	0	100	100
1	D	236/238 (99%)	230 (98%)	6 (2%)	0	100	100
2	C	207/215 (96%)	202 (98%)	5 (2%)	0	100	100
3	B	210/212 (99%)	204 (97%)	6 (3%)	0	100	100
3	E	212/212 (100%)	204 (96%)	7 (3%)	1 (0%)	24	15
4	F	207/216 (96%)	203 (98%)	4 (2%)	0	100	100
All	All	1310/1331 (98%)	1274 (97%)	35 (3%)	1 (0%)	48	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	151	ASN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	201/201 (100%)	201 (100%)	0	100	100
1	D	199/201 (99%)	199 (100%)	0	100	100
2	C	179/183 (98%)	179 (100%)	0	100	100
3	B	184/184 (100%)	184 (100%)	0	100	100
3	E	185/184 (100%)	185 (100%)	0	100	100
4	F	178/184 (97%)	178 (100%)	0	100	100
All	All	1126/1137 (99%)	1126 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	5	GLN
3	B	146	GLN
1	D	98	ASN
1	D	177	ASN
4	F	39	GLN
3	E	37	GLN
3	E	52	ASN
3	E	159	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	238/238 (100%)	0.67	24 (10%)	12 14	19, 42, 67, 74	2 (0%)
1	D	238/238 (100%)	0.68	11 (4%)	37 42	31, 44, 71, 79	0
2	C	209/215 (97%)	0.41	13 (6%)	26 30	22, 37, 67, 82	2 (0%)
3	B	211/212 (99%)	0.28	13 (6%)	26 30	24, 37, 65, 72	1 (0%)
3	E	212/212 (100%)	0.27	1 (0%)	87 91	26, 41, 54, 70	2 (0%)
4	F	210/216 (97%)	0.07	4 (1%)	66 72	22, 35, 52, 65	1 (0%)
All	All	1318/1331 (99%)	0.41	66 (5%)	34 38	19, 39, 66, 82	8 (0%)

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	105	ALA	4.4
2	C	128	PRO	4.1
2	C	189	SER	3.9
2	C	196	TYR	3.9
1	D	121	PHE	3.8
2	C	135	GLY	3.7
2	C	160	ALA	3.6
1	A	121	PHE	3.5
1	A	179	ILE	3.4
1	D	108	THR	3.3
4	F	129	SER	3.2
3	B	149	VAL	3.2
1	D	270	PRO	3.1
2	C	158	SER	2.9
2	C	193	THR	2.9
1	A	222	LEU	2.9
1	A	161	ARG	2.8
3	B	183	ALA	2.8
1	A	272	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	184	THR	2.8
1	A	229	VAL	2.8
4	F	42	GLY	2.7
2	C	215	PRO	2.7
2	C	136	GLY	2.7
1	A	232	PHE	2.6
1	D	122	THR	2.6
3	B	207	SER	2.6
1	D	228	LEU	2.6
2	C	195	THR	2.6
3	B	153	LEU	2.6
4	F	1	GLN	2.6
3	E	105	LEU	2.5
1	A	228	LEU	2.5
3	B	152	ALA	2.5
1	A	249	VAL	2.5
1	A	107	GLU	2.4
1	A	190	TYR	2.4
1	A	244	LEU	2.4
1	A	182	PHE	2.4
1	A	183	PHE	2.4
1	A	176	LEU	2.4
3	B	124	LEU	2.4
1	A	243	LEU	2.3
3	B	208	PHE	2.3
2	C	127	ALA	2.3
1	D	107	GLU	2.2
3	B	190	VAL	2.2
2	C	208	LYS	2.2
1	A	273	THR	2.2
3	B	180	LEU	2.2
1	D	120	GLY	2.2
2	C	148	PHE	2.2
1	A	175	LYS	2.2
3	B	188	HIS	2.1
1	D	182	PHE	2.1
1	A	269	LEU	2.1
3	B	192	ALA	2.1
3	B	193	CYS	2.1
4	F	193	THR	2.0
1	A	177	ASN	2.0
1	A	119	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	227	ASP	2.0
1	A	226	SER	2.0
3	B	181	SER	2.0
1	A	122	THR	2.0
1	D	181	GLY	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.