



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:29 AM GMT

PDB ID : 2X8R
Title : The structure of a family GH25 lysozyme from *Aspergillus fumigatus*
Authors : Korczynska, J.E.; Danielsen, S.; Schagerlof, U.; Turkenburg, J.P.; Davies, G.J.;
Wilson, K.S.; Taylor, E.J.
Deposited on : 2010-03-11
Resolution : 1.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

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

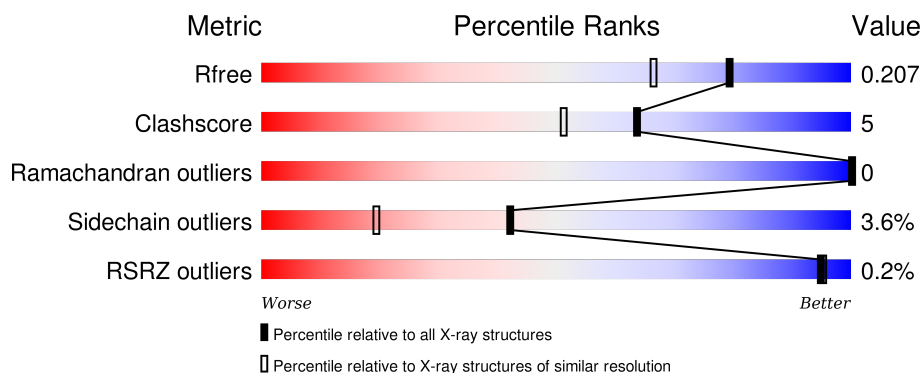
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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	210	 89% 10%
1	B	210	 86% 13%
1	C	210	 87% 12%
1	D	210	 90% 9%
1	E	210	 87% 11%

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Mol	Chain	Length	Quality of chain
1	F	210	 91% 6% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11208 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 GLYCOSYL HYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	9	0
			1693	1069	293	322	9			
1	B	210	Total	C	N	O	S	0	8	0
			1702	1073	298	322	9			
1	C	210	Total	C	N	O	S	0	10	0
			1720	1089	299	323	9			
1	D	209	Total	C	N	O	S	0	3	0
			1654	1047	287	312	8			
1	E	210	Total	C	N	O	S	0	3	0
			1653	1046	286	313	8			
1	F	209	Total	C	N	O	S	0	5	0
			1651	1047	285	310	9			

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Cl	0	0
			1	1		
2	E	1	Total	Cl	0	0
			1	1		
2	B	1	Total	Cl	0	0
			1	1		
2	C	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		
2	F	1	Total	Cl	0	0
			1	1		


- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	218	Total 218	O 218	0	0
3	B	221	Total 221	O 221	0	0
3	C	197	Total 197	O 197	0	0
3	D	171	Total 171	O 171	0	0
3	E	169	Total 169	O 169	0	0
3	F	153	Total 153	O 153	0	0

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: GLYCOSYL HYDROLASE

Chain A: 



• Molecule 1: GLYCOSYL HYDROLASE

Chain B: 



• Molecule 1: GLYCOSYL HYDROLASE

Chain C: 



• Molecule 1: GLYCOSYL HYDROLASE

Chain D: 



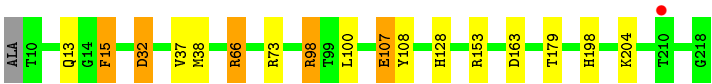
• Molecule 1: GLYCOSYL HYDROLASE

Chain E: 



• Molecule 1: GLYCOSYL HYDROLASE

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.65Å 111.72Å 119.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	81.50 – 1.70 40.75 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (81.50-1.70) 99.9 (40.75-1.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.6.0056	Depositor
R, R_{free}	0.169 , 0.207 0.169 , 0.207	Depositor DCC
R_{free} test set	5938 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	13.6	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	1 of 118583 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11208	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.30 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 7.2248e-03.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CL

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	0.84	0/1749	0.77	0/2376
1	B	0.92	1/1757 (0.1%)	0.82	1/2384 (0.0%)
1	C	0.87	0/1782	0.78	2/2416 (0.1%)
1	D	0.84	1/1709 (0.1%)	0.77	0/2321
1	E	0.87	2/1711 (0.1%)	0.72	0/2323
1	F	0.77	1/1715 (0.1%)	0.71	1/2331 (0.0%)
All	All	0.85	5/10423 (0.0%)	0.76	4/14151 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	107	GLU	CD-OE1	-5.55	1.19	1.25
1	E	139	SER	CB-OG	-5.51	1.35	1.42
1	D	96	ASP	C-O	-5.34	1.13	1.23
1	E	107	GLU	CD-OE1	-5.13	1.20	1.25
1	B	107	GLU	CD-OE2	-5.03	1.20	1.25

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	169	LEU	CB-CG-CD2	5.78	120.82	111.00
1	C	169	LEU	CB-CG-CD2	5.56	120.45	111.00
1	F	98	ARG	NE-CZ-NH1	5.21	122.90	120.30
1	C	98	ARG	NE-CZ-NH1	5.03	122.81	120.30

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	1693	0	1562	19	0
1	B	1702	0	1587	21	0
1	C	1720	0	1616	20	0
1	D	1654	0	1544	17	0
1	E	1653	0	1545	18	0
1	F	1651	0	1534	11	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	218	0	0	8	0
3	B	221	0	0	5	0
3	C	197	0	0	5	0
3	D	171	0	0	5	0
3	E	169	0	0	5	0
3	F	153	0	0	4	0
All	All	11208	0	9388	101	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 5.

All (101) 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:A:2111:HOH:O	1:B:98[B]:ARG:HD3	1.62	1.00
1:A:136:HIS:HE1	1:C:120:HIS:H	1.09	0.93
1:C:38[B]:MET:SD	1:C:100:LEU:HD12	2.13	0.89
1:A:136:HIS:CE1	1:C:120:HIS:H	1.96	0.84
1:C:88:LYS:HE3	3:C:2085:HOH:O	1.79	0.82
1:F:73[A]:ARG:HD2	3:F:2055:HOH:O	1.80	0.82
3:E:2140:HOH:O	1:F:98:ARG:HD2	1.84	0.76
1:D:84:LYS:HE3	1:E:11:THR:O	1.86	0.76
1:B:173[A]:SER:OG	1:B:174:SER:N	2.17	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38[B]:MET:SD	1:C:100:LEU:CD1	2.77	0.72
1:D:131[A]:VAL:HG22	1:D:142:PRO:HG2	1.71	0.71
1:A:73:ARG:HD2	3:A:2073:HOH:O	1.91	0.71
1:A:198:HIS:HD2	3:A:2081:HOH:O	1.72	0.71
1:E:80:SER:HB3	3:E:2069:HOH:O	1.91	0.70
1:B:88:LYS:HE2	3:B:2089:HOH:O	1.91	0.69
1:B:38[B]:MET:SD	1:B:100:LEU:HD12	2.33	0.68
1:B:135:HIS:HE1	3:B:2150:HOH:O	1.76	0.68
1:B:45[B]:THR:HG22	1:B:77:SER:OG	1.96	0.65
1:A:211:GLN:HE22	1:A:214[A]:LYS:HE2	1.62	0.64
1:C:161:PHE:HA	1:C:164[A]:LYS:HE2	1.78	0.64
1:A:38[B]:MET:SD	1:A:100:LEU:HD12	2.39	0.63
1:C:164[A]:LYS:H	1:C:164[A]:LYS:HZ3	1.45	0.62
1:F:128:HIS:HD2	3:F:2095:HOH:O	1.82	0.62
1:D:135:HIS:HE1	3:D:2111:HOH:O	1.83	0.61
1:B:32:ASP:OD2	1:B:198:HIS:HE1	1.84	0.60
1:E:198:HIS:HD2	3:E:2007:HOH:O	1.84	0.60
1:D:131[B]:VAL:HG13	1:D:142:PRO:HD2	1.84	0.59
1:A:32:ASP:OD2	1:A:198:HIS:HE1	1.87	0.58
1:E:114:THR:HG22	3:E:2091:HOH:O	2.02	0.58
1:B:38[B]:MET:SD	1:B:100:LEU:CD1	2.91	0.58
1:B:45[B]:THR:OG1	1:B:46:THR:HG23	2.04	0.58
1:D:114:THR:HB	3:D:2095:HOH:O	2.04	0.57
1:F:32:ASP:OD2	1:F:198:HIS:HE1	1.86	0.57
1:A:204:LYS:HE3	3:A:2006:HOH:O	2.04	0.57
1:F:38[B]:MET:SD	1:F:100:LEU:HD12	2.45	0.56
1:C:55:HIS:HD2	3:C:2019:HOH:O	1.88	0.56
1:B:45[A]:THR:HG22	3:B:2038:HOH:O	2.07	0.55
1:B:60:THR:HA	1:B:98[A]:ARG:NH2	2.22	0.54
1:B:32:ASP:OD2	1:B:198:HIS:CE1	2.59	0.54
1:A:25:ASN:ND2	1:A:28:ALA:H	2.06	0.54
1:D:32:ASP:OD2	1:D:198:HIS:HE1	1.92	0.53
1:D:10:THR:HG23	3:D:2029:HOH:O	2.09	0.53
1:D:13:GLN:NE2	1:D:204:LYS:NZ	2.57	0.53
1:E:45:THR:OG1	1:E:46:THR:HG23	2.09	0.53
1:E:138:THR:O	1:E:139:SER:CB	2.54	0.52
1:C:106:ILE:HG23	1:C:106:ILE:O	2.11	0.51
1:A:128[B]:HIS:CE1	1:A:132:ASN:HD21	2.28	0.51
1:E:32:ASP:OD2	1:E:198:HIS:HE1	1.94	0.51
1:D:84:LYS:CE	1:E:11:THR:O	2.56	0.51
1:A:211:GLN:NE2	1:A:214[A]:LYS:HE2	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:138:THR:O	1:D:139[B]:SER:HB2	2.11	0.50
1:A:128[A]:HIS:HE1	3:A:2173:HOH:O	1.94	0.50
1:B:55:HIS:HD2	3:B:2014:HOH:O	1.95	0.50
1:A:55:HIS:HD2	3:A:2021:HOH:O	1.94	0.50
1:D:55:HIS:HD2	3:D:2016:HOH:O	1.95	0.49
1:B:141:TRP:NE1	1:B:164:LYS:HD3	2.28	0.49
1:F:128:HIS:HE1	3:F:2118:HOH:O	1.95	0.48
1:E:55:HIS:HD2	3:E:2015:HOH:O	1.97	0.48
1:D:84:LYS:HZ1	1:E:11:THR:H	1.61	0.48
1:A:189[A]:ILE:HD13	1:A:205:PHE:HA	1.95	0.48
1:C:30:LYS:HE2	1:C:63:GLY:O	2.14	0.48
1:B:88:LYS:NZ	3:B:2092:HOH:O	2.46	0.47
1:B:9:ALA:O	1:B:208:PRO:HB3	2.14	0.47
1:C:136:HIS:HD2	3:C:2125:HOH:O	1.96	0.47
1:B:45[A]:THR:HG23	1:B:46:THR:HG23	1.95	0.47
1:E:72:ALA:HB2	1:E:104:LEU:HD11	1.97	0.47
1:A:159:LYS:HE3	3:A:2169:HOH:O	2.13	0.47
1:A:128[A]:HIS:HD2	3:A:2142:HOH:O	1.98	0.46
1:E:163:ASP:OD1	1:E:164:LYS:HG3	2.15	0.46
1:F:37:VAL:O	1:F:66:ARG:HA	2.16	0.46
1:C:32:ASP:OD2	1:C:198:HIS:HE1	1.99	0.46
1:C:131:VAL:HG13	1:C:142:PRO:HD2	1.98	0.45
1:F:15:PHE:O	1:F:37:VAL:HA	2.16	0.45
1:B:60:THR:HA	1:B:98[A]:ARG:HH22	1.81	0.45
1:D:131[A]:VAL:HG22	1:D:142:PRO:CG	2.43	0.45
1:A:38[B]:MET:SD	1:A:100:LEU:CD1	3.04	0.45
1:C:13:GLN:HE22	1:C:204:LYS:HE2	1.81	0.44
1:D:32:ASP:OD2	1:D:198:HIS:CE1	2.71	0.44
1:C:135:HIS:HE1	3:C:2127:HOH:O	2.01	0.44
1:B:131:VAL:HG22	1:B:142:PRO:HG2	2.01	0.43
1:C:47[B]:TYR:CE2	1:C:49:ASP:HA	2.54	0.43
1:E:40:LYS:HB2	1:E:69:TYR:CZ	2.55	0.42
1:C:88:LYS:HE2	3:C:2082:HOH:O	2.18	0.42
1:C:13:GLN:NE2	1:C:204:LYS:HE2	2.34	0.42
1:B:167:LEU:HD21	1:B:169:LEU:HG	2.01	0.42
1:A:168:VAL:HG13	1:A:189[A]:ILE:HB	2.02	0.42
1:F:13:GLN:NE2	1:F:204:LYS:HE2	2.35	0.42
1:B:162:GLY:HA2	1:B:184:TRP:CD1	2.55	0.42
1:F:107:GLU:HB2	1:F:108:TYR:H	1.71	0.42
1:E:131:VAL:HG22	1:E:142:PRO:HG2	2.00	0.41
1:E:25:ASN:ND2	1:E:28:ALA:H	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:15:PHE:O	1:D:37:VAL:HA	2.20	0.41
1:D:57:THR:HG22	1:D:61:LYS:HD2	2.02	0.41
1:C:38[B]:MET:SD	1:C:100:LEU:HD13	2.58	0.41
1:E:45:THR:CG2	1:E:77:SER:HB3	2.50	0.41
1:C:37:VAL:O	1:C:66:ARG:HA	2.21	0.41
1:F:153:ARG:NE	3:F:2112:HOH:O	2.51	0.41
1:E:138:THR:O	1:E:139:SER:HB3	2.19	0.41
1:E:74:PRO:HD2	1:E:116:TYR:CG	2.56	0.41
1:D:25:ASN:ND2	3:D:2017:HOH:O	2.27	0.40
1:A:25:ASN:HD22	1:A:25:ASN:C	2.25	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	216/210 (103%)	212 (98%)	4 (2%)	0	100	100
1	B	216/210 (103%)	213 (99%)	3 (1%)	0	100	100
1	C	218/210 (104%)	214 (98%)	4 (2%)	0	100	100
1	D	210/210 (100%)	203 (97%)	7 (3%)	0	100	100
1	E	211/210 (100%)	204 (97%)	7 (3%)	0	100	100
1	F	212/210 (101%)	208 (98%)	4 (2%)	0	100	100
All	All	1283/1260 (102%)	1254 (98%)	29 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	176/171 (103%)	169 (96%)	7 (4%)	38	16
1	B	177/171 (104%)	170 (96%)	7 (4%)	38	16
1	C	179/171 (105%)	170 (95%)	9 (5%)	30	11
1	D	172/171 (101%)	166 (96%)	6 (4%)	43	20
1	E	172/171 (101%)	165 (96%)	7 (4%)	37	15
1	F	170/171 (99%)	165 (97%)	5 (3%)	50	27
All	All	1046/1026 (102%)	1005 (96%)	41 (4%)	42	16

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	PHE
1	A	25	ASN
1	A	27	GLU
1	A	31	LYS
1	A	66	ARG
1	A	114[A]	THR
1	A	114[B]	THR
1	B	15	PHE
1	B	66	ARG
1	B	107	GLU
1	B	114[A]	THR
1	B	114[B]	THR
1	B	169	LEU
1	B	175	SER
1	C	15	PHE
1	C	25	ASN
1	C	66	ARG
1	C	114[A]	THR
1	C	114[B]	THR
1	C	164[A]	LYS
1	C	164[B]	LYS
1	C	169	LEU

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Mol	Chain	Res	Type
1	C	179	THR
1	D	15	PHE
1	D	66	ARG
1	D	107	GLU
1	D	114	THR
1	D	131[A]	VAL
1	D	131[B]	VAL
1	E	15	PHE
1	E	25	ASN
1	E	66	ARG
1	E	76	LYS
1	E	80	SER
1	E	114	THR
1	E	179	THR
1	F	15	PHE
1	F	32	ASP
1	F	66	ARG
1	F	163	ASP
1	F	179	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	25	ASN
1	A	35	GLN
1	A	55	HIS
1	A	132	ASN
1	A	136	HIS
1	A	198	HIS
1	B	13	GLN
1	B	55	HIS
1	B	135	HIS
1	B	198	HIS
1	B	211	GLN
1	C	13	GLN
1	C	25	ASN
1	C	55	HIS
1	C	135	HIS
1	C	136	HIS
1	C	198	HIS
1	D	13	GLN

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Mol	Chain	Res	Type
1	D	20	HIS
1	D	55	HIS
1	D	135	HIS
1	D	198	HIS
1	E	13	GLN
1	E	25	ASN
1	E	55	HIS
1	E	198	HIS
1	F	13	GLN
1	F	55	HIS
1	F	128	HIS
1	F	198	HIS
1	F	211	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	209/210 (99%)	-0.37	0 100 100	5, 10, 17, 25	0
1	B	210/210 (100%)	-0.28	2 (0%) 84 87	5, 10, 20, 33	0
1	C	210/210 (100%)	-0.21	0 100 100	7, 13, 22, 29	0
1	D	209/210 (99%)	-0.27	0 100 100	7, 13, 25, 29	0
1	E	210/210 (100%)	-0.26	0 100 100	7, 13, 23, 28	0
1	F	209/210 (99%)	-0.00	1 (0%) 91 93	9, 17, 30, 38	0
All	All	1257/1260 (99%)	-0.23	3 (0%) 95 95	5, 12, 24, 38	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	9	ALA	3.7
1	B	174	SER	2.6
1	F	210	THR	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CL	A	1219	1/1	1.00	0.09	0.75	9,9,9,9	0
2	CL	B	1219	1/1	1.00	0.07	-0.45	9,9,9,9	0
2	CL	F	1219	1/1	1.00	0.06	-1.18	10,10,10,10	0
2	CL	D	1219	1/1	0.99	0.07	-1.22	13,13,13,13	0
2	CL	C	1219	1/1	1.00	0.05	-1.68	9,9,9,9	0
2	CL	E	1219	1/1	0.99	0.04	-2.42	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.