



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:17 AM GMT

PDB ID : 3A6F
Title : W174F mutant creatininase, Type II
Authors : Nakajima, Y.; Yamashita, K.; Ito, K.; Yoshimoto, T.
Deposited on : 2009-08-31
Resolution : 1.78 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

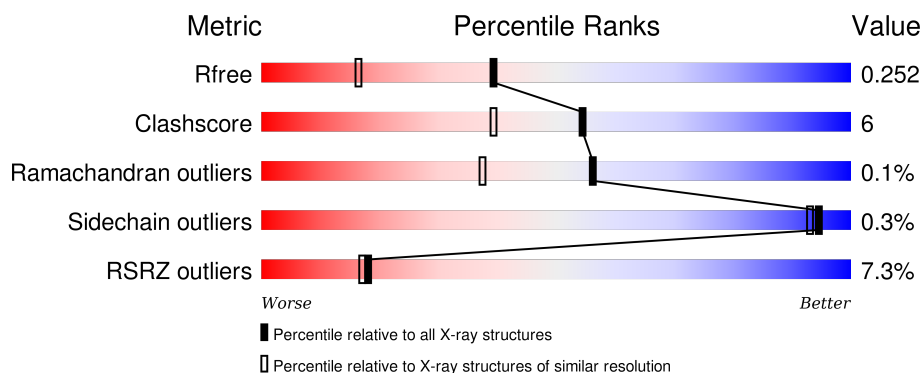
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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	6655 (1.80-1.76)
Clashscore	102246	7658 (1.80-1.76)
Ramachandran outliers	100387	7570 (1.80-1.76)
Sidechain outliers	100360	7569 (1.80-1.76)
RSRZ outliers	91569	6671 (1.80-1.76)

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	260	<div> <div>8%</div> <div>83%</div> <div>16%</div> </div>
1	B	260	<div> <div>9%</div> <div>85%</div> <div>14%</div> </div>
1	C	260	<div> <div>8%</div> <div>83%</div> <div>15%</div> </div>
1	D	260	<div> <div>8%</div> <div>84%</div> <div>15%</div> </div>
1	E	260	<div> <div>6%</div> <div>87%</div> <div>12%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	260	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a red segment at the beginning labeled '5%', a green segment in the middle labeled '85%', and a yellow segment at the end labeled '13%'. A small grey dot is located at the far right end of the bar.

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12656 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 Creatinine amidohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			1984	1271	337	365	11			
1	B	257	Total	C	N	O	S	0	0	0
			1981	1269	336	365	11			
1	C	257	Total	C	N	O	S	0	0	0
			1977	1267	335	364	11			
1	D	257	Total	C	N	O	S	0	0	0
			1977	1266	335	365	11			
1	E	257	Total	C	N	O	S	0	0	0
			1970	1263	334	362	11			
1	F	256	Total	C	N	O	S	0	0	0
			1961	1257	332	361	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	174	PHE	TRP	ENGINEERED	UNP P83772
B	174	PHE	TRP	ENGINEERED	UNP P83772
C	174	PHE	TRP	ENGINEERED	UNP P83772
D	174	PHE	TRP	ENGINEERED	UNP P83772
E	174	PHE	TRP	ENGINEERED	UNP P83772
F	174	PHE	TRP	ENGINEERED	UNP P83772

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Mn	0	0
			1	1		
2	E	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		

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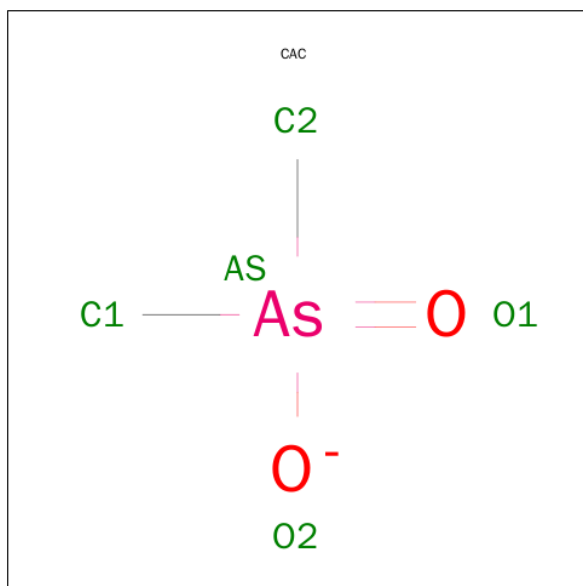
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total 1	Mn 1	0	0
2	A	1	Total 1	Mn 1	0	0
2	F	1	Total 1	Mn 1	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total 1	Zn 1	0	0
3	E	1	Total 1	Zn 1	0	0
3	B	1	Total 1	Zn 1	0	0
3	C	1	Total 1	Zn 1	0	0
3	A	1	Total 1	Zn 1	0	0
3	F	1	Total 1	Zn 1	0	0

- Molecule 4 is CACODYLATE ION (three-letter code: CAC) (formula: C₂H₆AsO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	As	C	O	0	0
			5	1	2	2		
4	B	1	Total	As	C	O	0	0
			5	1	2	2		
4	C	1	Total	As	C	O	0	0
			5	1	2	2		
4	D	1	Total	As	C	O	0	0
			5	1	2	2		
4	E	1	Total	As	C	O	0	0
			5	1	2	2		
4	F	1	Total	As	C	O	0	0
			5	1	2	2		

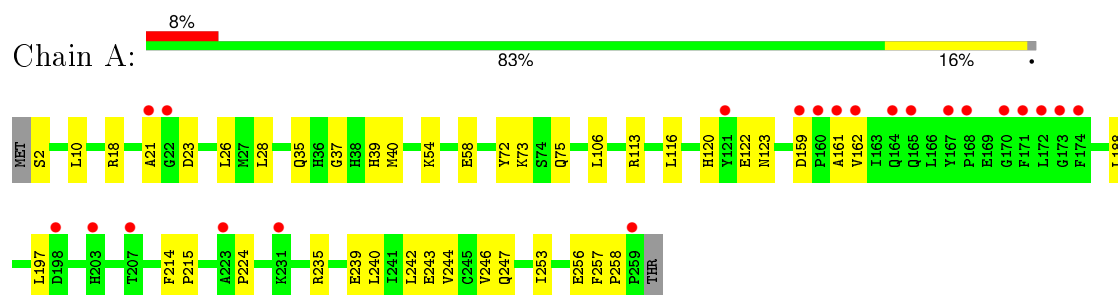
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	111	Total	O	0	0
			111	111		
5	B	124	Total	O	0	0
			124	124		
5	C	126	Total	O	0	0
			126	126		
5	D	128	Total	O	0	0
			128	128		
5	E	125	Total	O	0	0
			125	125		
5	F	150	Total	O	0	0
			150	150		

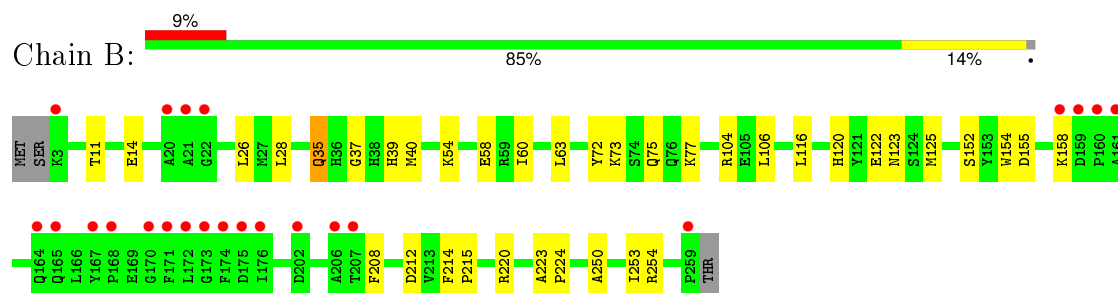
3 Residue-property plots

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.

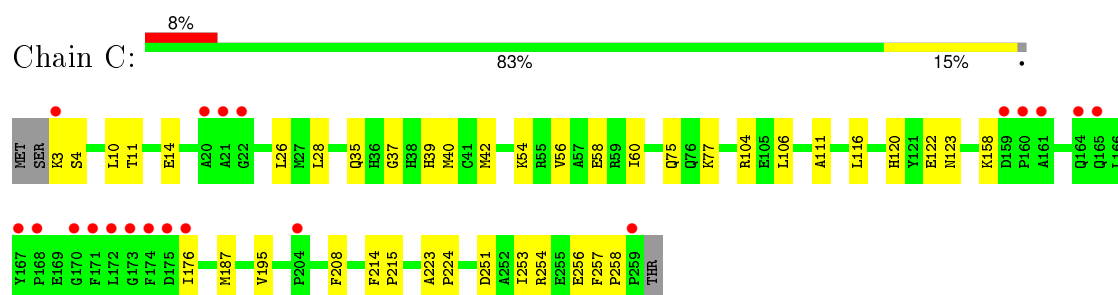
- Molecule 1: Creatinine amidohydrolase



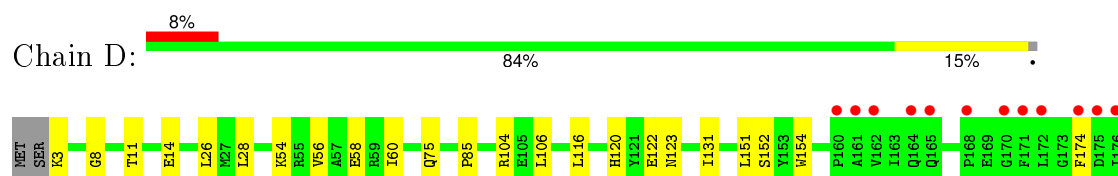
- Molecule 1: Creatinine amidohydrolase

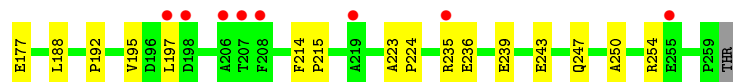


- Molecule 1: Creatinine amidohydrolase

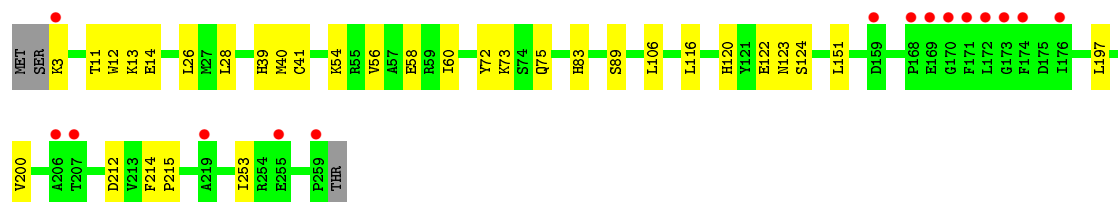
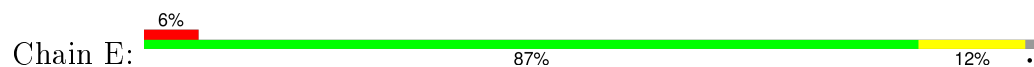


- Molecule 1: Creatinine amidohydrolase

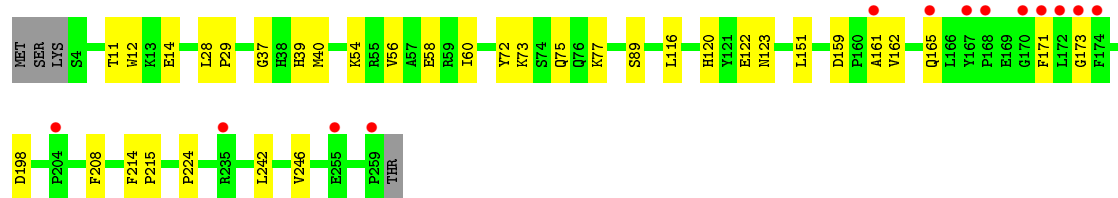
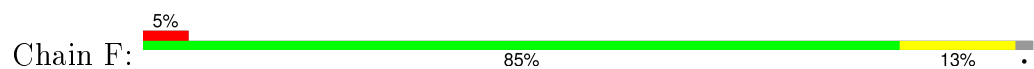




- Molecule 1: Creatinine amidohydrolase



- Molecule 1: Creatinine amidohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	164.30Å 164.30Å 163.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.78 32.59 – 1.78	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-1.78) 99.8 (32.59-1.78)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 1.78Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.235 , 0.257 0.232 , 0.252	Depositor DCC
R_{free} test set	12155 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	21.8	Xtriage
Anisotropy	0.153	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 59.8	EDS
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 242442 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12656	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% 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, 0.375 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: CAC, ZN, MN

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	0.32	0/2032	0.57	0/2764
1	B	0.32	0/2029	0.58	0/2760
1	C	0.33	0/2025	0.57	0/2755
1	D	0.33	0/2025	0.57	0/2756
1	E	0.34	0/2018	0.58	0/2747
1	F	0.33	0/2009	0.58	0/2736
All	All	0.33	0/12138	0.58	0/16518

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1984	0	1967	27	0
1	B	1981	0	1964	26	0
1	C	1977	0	1958	27	0
1	D	1977	0	1953	30	0
1	E	1970	0	1945	23	0
1	F	1961	0	1932	21	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	5	0	0	1	0
4	B	5	0	0	1	0
4	C	5	0	0	1	0
4	D	5	0	0	0	0
4	E	5	0	0	1	0
4	F	5	0	0	0	0
5	A	111	0	0	1	0
5	B	124	0	0	3	0
5	C	126	0	0	3	0
5	D	128	0	0	4	0
5	E	125	0	0	2	0
5	F	150	0	0	1	0
All	All	12656	0	11719	152	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 6.

All (152) 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:E:120:HIS:HB3	1:E:123:ASN:ND2	1.94	0.81
1:A:120:HIS:HB3	1:A:123:ASN:ND2	2.00	0.76
1:C:28:LEU:HD23	1:C:116:LEU:HD21	1.69	0.75
1:A:54:LYS:O	1:A:58:GLU:HG3	1.90	0.72
1:D:188:LEU:O	1:D:192:PRO:HG3	1.91	0.70
1:E:28:LEU:HD23	1:E:116:LEU:HD21	1.75	0.68
1:B:72:TYR:H	1:B:123:ASN:ND2	1.93	0.66
1:F:198:ASP:HB2	5:F:1336:HOH:O	1.96	0.65
1:C:54:LYS:O	1:C:58:GLU:HG3	1.97	0.65
1:C:120:HIS:HB3	1:C:123:ASN:ND2	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:28:LEU:HD23	1:F:116:LEU:HD21	1.78	0.65
1:D:120:HIS:HB3	1:D:123:ASN:ND2	2.13	0.64
1:E:3:LYS:HE2	5:E:1666:HOH:O	1.99	0.63
1:D:188:LEU:HD21	1:D:197:LEU:HD11	1.81	0.62
1:C:4:SER:HB2	5:C:1586:HOH:O	1.98	0.62
1:A:23:ASP:HB2	1:A:113:ARG:NH2	2.16	0.60
1:F:72:TYR:O	1:F:120:HIS:HE1	1.84	0.60
1:F:120:HIS:HB3	1:F:123:ASN:ND2	2.16	0.60
1:B:28:LEU:HD23	1:B:116:LEU:HD21	1.84	0.60
1:F:54:LYS:O	1:F:58:GLU:HG3	2.03	0.59
1:E:120:HIS:HA	4:E:303:CAC:O1	2.03	0.58
1:B:54:LYS:O	1:B:58:GLU:HG3	2.04	0.58
1:A:243:GLU:O	1:A:247:GLN:HG3	2.03	0.58
1:A:28:LEU:HD23	1:A:116:LEU:HD21	1.85	0.58
1:D:235:ARG:O	1:D:239:GLU:HG3	2.04	0.58
1:F:28:LEU:HD23	1:F:116:LEU:CD2	2.33	0.57
1:C:120:HIS:HA	4:C:303:CAC:O1	2.04	0.57
1:D:243:GLU:HG3	1:D:247:GLN:HE21	1.70	0.57
1:C:104:ARG:HD2	5:C:1454:HOH:O	2.05	0.56
1:B:11:THR:OG1	1:B:14:GLU:HG3	2.06	0.55
1:D:195:VAL:HG12	1:D:197:LEU:HD12	1.89	0.55
1:B:220:ARG:HD3	5:B:1610:HOH:O	2.07	0.54
1:F:161:ALA:O	1:F:165:GLN:HG2	2.07	0.54
1:F:56:VAL:O	1:F:60:ILE:HG13	2.07	0.54
1:C:26:LEU:HD13	1:C:106:LEU:HD22	1.91	0.53
1:C:56:VAL:O	1:C:60:ILE:HG12	2.08	0.52
1:D:11:THR:OG1	1:D:14:GLU:HG3	2.09	0.52
1:B:72:TYR:H	1:B:123:ASN:HD21	1.57	0.52
1:E:54:LYS:O	1:E:58:GLU:HG3	2.10	0.52
1:E:214:PHE:HA	1:E:215:PRO:C	2.30	0.52
1:F:151:LEU:HD12	1:F:151:LEU:C	2.30	0.52
1:E:12:TRP:CZ2	1:E:13:LYS:HD3	2.46	0.51
1:C:26:LEU:CD1	1:C:106:LEU:HD22	2.40	0.51
1:B:223:ALA:N	1:B:224:PRO:HD2	2.26	0.51
1:D:214:PHE:HA	1:D:215:PRO:C	2.32	0.51
1:E:56:VAL:O	1:E:60:ILE:HG12	2.11	0.51
1:D:3:LYS:N	5:D:1238:HOH:O	2.44	0.50
1:D:223:ALA:HB3	1:D:224:PRO:HD3	1.92	0.50
1:E:60:ILE:HD12	1:E:253:ILE:HG21	1.93	0.50
1:B:250:ALA:O	1:B:254:ARG:HG3	2.11	0.50
1:F:11:THR:OG1	1:F:14:GLU:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:256:GLU:C	1:C:258:PRO:HD3	2.32	0.50
1:D:54:LYS:O	1:D:58:GLU:HG3	2.12	0.50
1:D:116:LEU:HD12	1:D:131:ILE:HD11	1.94	0.50
1:D:188:LEU:HD21	1:D:197:LEU:CD1	2.42	0.49
1:B:75:GLN:HG2	1:B:122:GLU:HG2	1.92	0.49
1:F:159:ASP:OD2	1:F:162:VAL:HG23	2.13	0.49
1:C:214:PHE:HA	1:C:215:PRO:C	2.33	0.49
1:D:75:GLN:HG2	1:D:122:GLU:HG2	1.94	0.49
1:E:28:LEU:HD23	1:E:116:LEU:CD2	2.42	0.49
1:A:120:HIS:HA	4:A:303:CAC:O2	2.13	0.49
1:B:214:PHE:HA	1:B:215:PRO:C	2.32	0.49
1:F:39:HIS:CD2	1:F:40:MET:HG3	2.47	0.49
1:A:240:LEU:O	1:A:244:VAL:HG23	2.12	0.48
1:D:104:ARG:HD2	5:D:1681:HOH:O	2.13	0.48
1:D:26:LEU:HD13	1:D:106:LEU:HD22	1.95	0.48
1:A:256:GLU:C	1:A:258:PRO:HD3	2.33	0.48
1:A:214:PHE:HA	1:A:215:PRO:C	2.33	0.48
1:F:75:GLN:HG2	1:F:122:GLU:HG2	1.95	0.48
1:F:214:PHE:HA	1:F:215:PRO:C	2.33	0.48
1:B:35:GLN:NE2	1:B:37:GLY:H	2.11	0.48
1:A:39:HIS:CD2	1:A:40:MET:HG3	2.48	0.48
1:E:39:HIS:CD2	1:E:40:MET:HG3	2.49	0.47
1:C:251:ASP:HA	1:C:254:ARG:HH11	1.79	0.47
1:B:120:HIS:HA	4:B:303:CAC:O2	2.15	0.47
1:B:26:LEU:HD23	1:B:63:LEU:HB2	1.96	0.47
1:C:60:ILE:HD12	1:C:253:ILE:HG21	1.97	0.47
1:D:28:LEU:HD23	1:D:116:LEU:HD21	1.96	0.46
1:B:26:LEU:HD13	1:B:106:LEU:HD22	1.96	0.46
1:B:39:HIS:CD2	1:B:40:MET:HG3	2.49	0.46
1:E:151:LEU:C	1:E:151:LEU:HD12	2.34	0.46
1:A:159:ASP:HB3	1:A:162:VAL:HG23	1.98	0.46
1:B:120:HIS:HB3	1:B:123:ASN:OD1	2.16	0.46
1:E:120:HIS:HB3	1:E:123:ASN:HD21	1.77	0.46
1:B:60:ILE:HD12	1:B:253:ILE:HG21	1.98	0.45
1:E:11:THR:OG1	1:E:14:GLU:HG3	2.17	0.45
1:E:197:LEU:HA	1:E:200:VAL:HG23	1.99	0.45
1:A:26:LEU:CD1	1:A:106:LEU:HD22	2.46	0.45
1:D:104:ARG:HB3	5:D:1681:HOH:O	2.17	0.45
1:E:83:HIS:CD2	1:E:83:HIS:H	2.35	0.45
1:B:120:HIS:HD2	1:B:123:ASN:ND2	2.15	0.45
1:A:37:GLY:HA2	1:A:224:PRO:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:ARG:NH2	1:A:239:GLU:OE2	2.49	0.44
1:B:104:ARG:HD2	5:B:1365:HOH:O	2.17	0.44
1:A:2:SER:O	1:A:10:LEU:HD23	2.17	0.44
1:D:85:PRO:HD3	5:D:1600:HOH:O	2.16	0.44
1:A:159:ASP:OD2	1:A:162:VAL:HG23	2.17	0.44
1:F:73:LYS:HA	1:F:89:SER:CB	2.47	0.44
1:C:77:LYS:HE2	1:C:208:PHE:HB3	1.99	0.44
1:D:174:PHE:HD1	1:D:177:GLU:HG3	1.83	0.44
1:C:111:ALA:HB1	5:C:1488:HOH:O	2.18	0.44
1:A:72:TYR:CG	1:A:73:LYS:N	2.86	0.44
1:D:236:GLU:N	1:D:236:GLU:OE1	2.48	0.44
1:B:72:TYR:CG	1:B:73:LYS:N	2.85	0.44
1:A:28:LEU:HD23	1:A:116:LEU:CD2	2.48	0.44
1:C:75:GLN:HG2	1:C:122:GLU:HG2	2.00	0.43
1:C:223:ALA:HB3	1:C:224:PRO:HD3	1.99	0.43
1:C:176:ILE:C	1:C:176:ILE:HD12	2.37	0.43
1:A:18:ARG:O	1:A:21:ALA:HB3	2.18	0.43
1:B:104:ARG:HB3	5:B:1365:HOH:O	2.18	0.43
1:B:155:ASP:O	1:B:158:LYS:HE3	2.18	0.43
1:E:83:HIS:HE1	5:E:1109:HOH:O	2.00	0.43
1:A:242:LEU:O	1:A:246:VAL:HG23	2.18	0.43
1:A:253:ILE:HG23	1:A:257:PHE:HD2	1.83	0.43
1:E:73:LYS:HA	1:E:89:SER:CB	2.49	0.43
1:C:60:ILE:CD1	1:C:253:ILE:HG21	2.49	0.43
1:A:235:ARG:HG2	1:A:239:GLU:OE2	2.18	0.43
1:B:152:SER:HB3	1:B:154:TRP:CE2	2.53	0.43
1:A:159:ASP:HB3	1:A:162:VAL:CG2	2.49	0.42
1:A:35:GLN:HB2	1:F:12:TRP:CD1	2.54	0.42
1:E:60:ILE:CD1	1:E:253:ILE:HG21	2.49	0.42
1:D:56:VAL:O	1:D:60:ILE:HG12	2.19	0.42
1:D:152:SER:HB3	1:D:154:TRP:CE2	2.54	0.42
1:A:159:ASP:OD2	1:A:161:ALA:HB3	2.19	0.42
1:D:151:LEU:C	1:D:151:LEU:HD12	2.40	0.42
1:B:77:LYS:HE2	1:B:208:PHE:CB	2.49	0.42
1:C:35:GLN:HA	1:C:42:MET:HG2	2.02	0.42
1:C:3:LYS:HG3	1:C:10:LEU:HD22	2.02	0.42
1:D:26:LEU:CD1	1:D:106:LEU:HD22	2.49	0.42
1:D:28:LEU:HD23	1:D:116:LEU:CD2	2.49	0.42
1:F:242:LEU:O	1:F:246:VAL:HG23	2.20	0.41
1:C:11:THR:OG1	1:C:14:GLU:HG3	2.21	0.41
1:B:28:LEU:HD23	1:B:116:LEU:CD2	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:HIS:CD2	1:C:40:MET:HG3	2.55	0.41
1:A:75:GLN:HG2	1:A:122:GLU:HG2	2.03	0.41
1:E:26:LEU:CD1	1:E:106:LEU:HD22	2.50	0.41
1:E:72:TYR:CG	1:E:73:LYS:N	2.89	0.41
1:D:243:GLU:CG	1:D:247:GLN:HE21	2.34	0.41
1:C:187:MET:HB3	1:C:195:VAL:CG2	2.51	0.41
1:F:77:LYS:HE2	1:F:208:PHE:HB3	2.02	0.40
1:C:28:LEU:HD23	1:C:116:LEU:CD2	2.47	0.40
1:D:188:LEU:HD11	1:D:197:LEU:HD11	2.02	0.40
1:C:37:GLY:HA2	1:C:224:PRO:O	2.21	0.40
1:D:250:ALA:O	1:D:254:ARG:HG3	2.20	0.40
1:D:8:GLY:O	1:E:41:CYS:HB2	2.21	0.40
1:F:28:LEU:HA	1:F:29:PRO:HD3	1.91	0.40
5:A:1120:HOH:O	1:B:125:MET:HG3	2.20	0.40
1:A:188:LEU:HD11	1:A:197:LEU:HD11	2.04	0.40
1:C:257:PHE:N	1:C:258:PRO:HD3	2.37	0.40
1:F:171:PHE:CZ	1:F:173:GLY:HA2	2.56	0.40
1:F:37:GLY:HA2	1:F:224:PRO:O	2.22	0.40
1:E:75:GLN:HG2	1:E:122:GLU:HG2	2.02	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	256/260 (98%)	247 (96%)	9 (4%)	0	100	100
1	B	255/260 (98%)	249 (98%)	6 (2%)	0	100	100
1	C	255/260 (98%)	248 (97%)	7 (3%)	0	100	100
1	D	255/260 (98%)	248 (97%)	7 (3%)	0	100	100
1	E	255/260 (98%)	248 (97%)	6 (2%)	1 (0%)	39	22

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	254/260 (98%)	248 (98%)	6 (2%)	0	100	100
All	All	1530/1560 (98%)	1488 (97%)	41 (3%)	1 (0%)	56	37

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	124	SER

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	211/215 (98%)	211 (100%)	0	100	100
1	B	211/215 (98%)	209 (99%)	2 (1%)	84	78
1	C	210/215 (98%)	209 (100%)	1 (0%)	92	89
1	D	210/215 (98%)	210 (100%)	0	100	100
1	E	208/215 (97%)	207 (100%)	1 (0%)	92	89
1	F	207/215 (96%)	207 (100%)	0	100	100
All	All	1257/1290 (97%)	1253 (100%)	4 (0%)	94	93

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	35	GLN
1	B	212	ASP
1	C	158	LYS
1	E	212	ASP

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

Mol	Chain	Res	Type
1	A	69	GLN

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Mol	Chain	Res	Type
1	A	83	HIS
1	A	165	GLN
1	B	35	GLN
1	B	123	ASN
1	D	247	GLN
1	E	83	HIS
1	F	120	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 12 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	CAC	A	303	3,2	0,4,4	0.00	-	0,6,6	0.00	-
4	CAC	B	303	3,2	0,4,4	0.00	-	0,6,6	0.00	-
4	CAC	C	303	3,2	0,4,4	0.00	-	0,6,6	0.00	-
4	CAC	D	303	3,2	0,4,4	0.00	-	0,6,6	0.00	-
4	CAC	E	303	3,2	0,4,4	0.00	-	0,6,6	0.00	-
4	CAC	F	303	3,2	0,4,4	0.00	-	0,6,6	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	CAC	A	303	3,2	-	0/0/0/0	0/0/0/0
4	CAC	B	303	3,2	-	0/0/0/0	0/0/0/0
4	CAC	C	303	3,2	-	0/0/0/0	0/0/0/0
4	CAC	D	303	3,2	-	0/0/0/0	0/0/0/0
4	CAC	E	303	3,2	-	0/0/0/0	0/0/0/0
4	CAC	F	303	3,2	-	0/0/0/0	0/0/0/0

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.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	303	CAC	1	0
4	B	303	CAC	1	0
4	C	303	CAC	1	0
4	E	303	CAC	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	258/260 (99%)	0.46	22 (8%)	13 12	12, 24, 47, 59	0
1	B	257/260 (98%)	0.38	23 (8%)	12 11	14, 22, 43, 56	0
1	C	257/260 (98%)	0.45	20 (7%)	16 15	13, 23, 44, 54	0
1	D	257/260 (98%)	0.35	20 (7%)	16 15	12, 23, 43, 53	0
1	E	257/260 (98%)	0.39	15 (5%)	26 24	11, 21, 41, 56	0
1	F	256/260 (98%)	0.33	13 (5%)	32 30	12, 20, 41, 58	0
All	All	1542/1560 (98%)	0.39	113 (7%)	18 17	11, 22, 44, 59	0

All (113) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	172	LEU	9.1
1	F	172	LEU	9.0
1	F	171	PHE	8.6
1	A	171	PHE	8.3
1	D	171	PHE	7.9
1	E	172	LEU	7.7
1	B	171	PHE	7.7
1	E	171	PHE	7.5
1	C	171	PHE	7.2
1	C	172	LEU	6.7
1	F	170	GLY	5.9
1	C	22	GLY	4.7
1	F	168	PRO	4.4
1	D	176	ILE	4.4
1	B	259	PRO	4.3
1	C	259	PRO	4.3
1	E	173	GLY	4.3
1	C	160	PRO	4.2
1	A	170	GLY	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	173	GLY	3.9
1	B	172	LEU	3.9
1	A	165	GLN	3.9
1	A	21	ALA	3.8
1	C	167	TYR	3.8
1	C	176	ILE	3.8
1	D	172	LEU	3.7
1	C	173	GLY	3.7
1	A	164	GLN	3.7
1	A	168	PRO	3.7
1	D	168	PRO	3.7
1	B	175	ASP	3.7
1	A	161	ALA	3.6
1	B	174	PHE	3.6
1	E	176	ILE	3.5
1	E	206	ALA	3.5
1	D	197	LEU	3.4
1	D	207	THR	3.4
1	D	164	GLN	3.3
1	C	21	ALA	3.3
1	F	161	ALA	3.3
1	F	173	GLY	3.3
1	B	170	GLY	3.2
1	D	206	ALA	3.2
1	B	164	GLN	3.2
1	A	160	PRO	3.1
1	E	174	PHE	3.1
1	B	22	GLY	3.1
1	B	20	ALA	3.1
1	F	165	GLN	3.1
1	D	174	PHE	3.0
1	C	161	ALA	3.0
1	F	174	PHE	3.0
1	C	168	PRO	2.9
1	B	207	THR	2.9
1	A	231	LYS	2.9
1	A	203	HIS	2.9
1	B	176	ILE	2.9
1	E	170	GLY	2.8
1	C	175	ASP	2.8
1	E	207	THR	2.8
1	A	223	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	158	LYS	2.7
1	A	207	THR	2.7
1	D	175	ASP	2.7
1	B	173	GLY	2.7
1	C	159	ASP	2.6
1	A	174	PHE	2.6
1	B	21	ALA	2.6
1	C	174	PHE	2.6
1	D	170	GLY	2.6
1	B	160	PRO	2.6
1	D	198	ASP	2.5
1	E	219	ALA	2.5
1	A	159	ASP	2.5
1	B	159	ASP	2.5
1	D	235	ARG	2.4
1	E	3	LYS	2.4
1	A	121	TYR	2.4
1	F	167	TYR	2.4
1	C	3	LYS	2.4
1	A	259	PRO	2.4
1	F	204	PRO	2.4
1	B	206	ALA	2.4
1	A	167	TYR	2.4
1	B	161	ALA	2.3
1	E	169	GLU	2.3
1	E	259	PRO	2.3
1	E	255	GLU	2.3
1	B	165	GLN	2.3
1	F	235	ARG	2.3
1	F	259	PRO	2.3
1	A	162	VAL	2.2
1	D	165	GLN	2.2
1	D	160	PRO	2.2
1	E	168	PRO	2.2
1	A	22	GLY	2.2
1	C	170	GLY	2.2
1	D	161	ALA	2.2
1	D	219	ALA	2.2
1	D	208	PHE	2.1
1	F	255	GLU	2.1
1	D	162	VAL	2.1
1	A	198	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	165	GLN	2.1
1	E	159	ASP	2.1
1	B	3	LYS	2.1
1	B	202	ASP	2.1
1	D	255	GLU	2.1
1	B	168	PRO	2.0
1	B	167	TYR	2.0
1	C	20	ALA	2.0
1	C	204	PRO	2.0
1	C	164	GLN	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 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(Å ²)	Q<0.9
4	CAC	B	303	5/5	0.98	0.13	1.80	30,30,35,35	0
4	CAC	A	303	5/5	0.96	0.14	1.54	38,38,40,40	0
4	CAC	E	303	5/5	0.97	0.12	1.48	34,34,36,37	0
4	CAC	C	303	5/5	0.98	0.11	0.10	35,37,38,39	0
2	MN	F	300	1/1	0.94	0.10	-0.05	36,36,36,36	0
4	CAC	F	303	5/5	0.98	0.09	-0.44	26,30,32,34	0
4	CAC	D	303	5/5	0.98	0.09	-0.52	32,33,35,36	0
2	MN	B	300	1/1	0.97	0.08	-0.62	39,39,39,39	0
2	MN	D	300	1/1	0.98	0.09	-0.85	40,40,40,40	0
2	MN	A	300	1/1	0.99	0.03	-2.80	34,34,34,34	0
2	MN	E	300	1/1	0.99	0.05	-3.14	33,33,33,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZN	F	301	1/1	1.00	0.04	-3.23	24,24,24,24	0
3	ZN	C	301	1/1	1.00	0.05	-3.31	25,25,25,25	0
3	ZN	B	301	1/1	0.99	0.04	-3.63	26,26,26,26	0
2	MN	C	300	1/1	0.99	0.04	-3.84	31,31,31,31	0
3	ZN	A	301	1/1	0.99	0.03	-4.39	27,27,27,27	0
3	ZN	E	301	1/1	0.99	0.02	-6.20	25,25,25,25	0
3	ZN	D	301	1/1	0.99	0.04	-7.29	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.