



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:17 AM GMT

PDB ID : 3A6H
Title : W154A mutant creatininase
Authors : Nakajima, Y.; Yamashita, K.; Ito, K.; Yoshimoto, T.
Deposited on : 2009-08-31
Resolution : 2.00 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

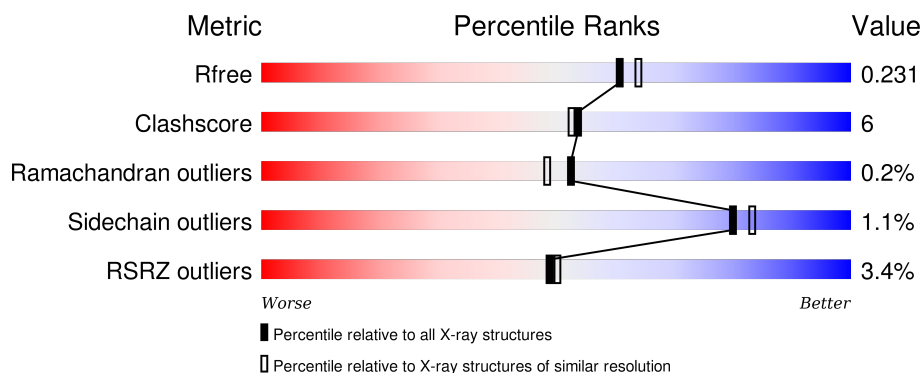
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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>90%</div> <div>8%</div> <div>.</div> </div>
1	B	260	<div> <div>2%</div> <div>89%</div> <div>9%</div> <div>.</div> </div>
1	C	260	<div> <div>2%</div> <div>83%</div> <div>15%</div> <div>..</div> </div>
1	D	260	<div> <div>2%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>
1	E	260	<div> <div>7%</div> <div>74%</div> <div>23%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	260	

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
4	CL	A	302	-	-	-	X
4	CL	D	305	-	-	-	X
4	CL	F	307	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12628 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	257	Total	C	N	O	S	0	0	0
			1979	1265	336	367	11			
1	B	256	Total	C	N	O	S	0	0	0
			1970	1259	334	366	11			
1	C	256	Total	C	N	O	S	0	0	0
			1966	1256	333	366	11			
1	D	257	Total	C	N	O	S	0	0	0
			1975	1262	335	367	11			
1	E	256	Total	C	N	O	S	0	0	0
			1960	1251	334	364	11			
1	F	257	Total	C	N	O	S	0	0	0
			1979	1265	336	367	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	154	ALA	TRP	ENGINEERED	UNP P83772
B	154	ALA	TRP	ENGINEERED	UNP P83772
C	154	ALA	TRP	ENGINEERED	UNP P83772
D	154	ALA	TRP	ENGINEERED	UNP P83772
E	154	ALA	TRP	ENGINEERED	UNP P83772
F	154	ALA	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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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 CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	157	Total 157	O 157	0	0
5	B	149	Total 149	O 149	0	0
5	C	127	Total 127	O 127	0	0
5	D	153	Total 153	O 153	0	0
5	E	86	Total 86	O 86	0	0
5	F	109	Total 109	O 109	0	0

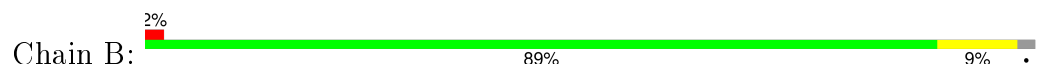
3 Residue-property plots [i](#)

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.

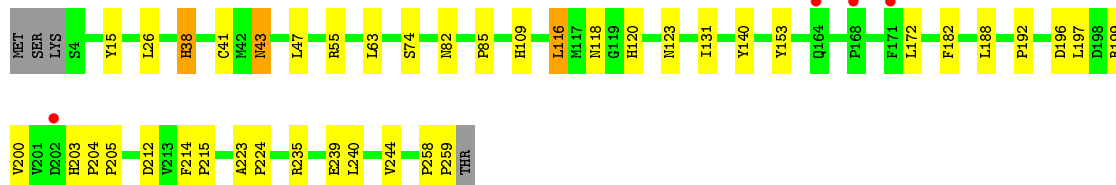
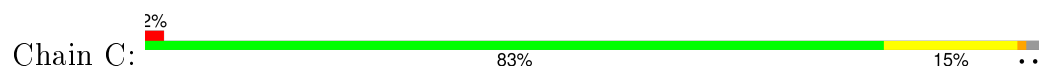
- Molecule 1: Creatinine amidohydrolase



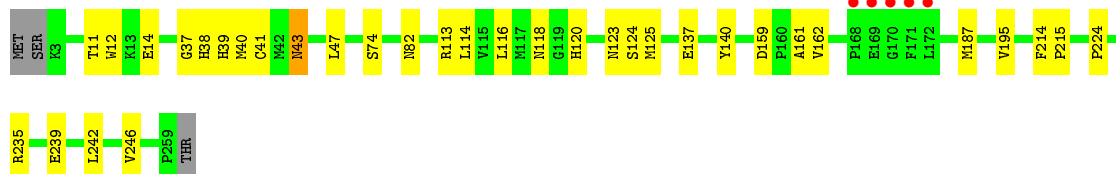
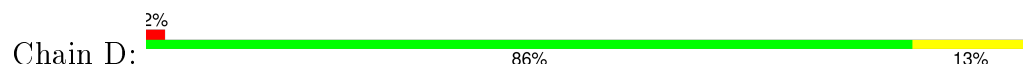
- Molecule 1: Creatinine amidohydrolase



- Molecule 1: Creatinine amidohydrolase

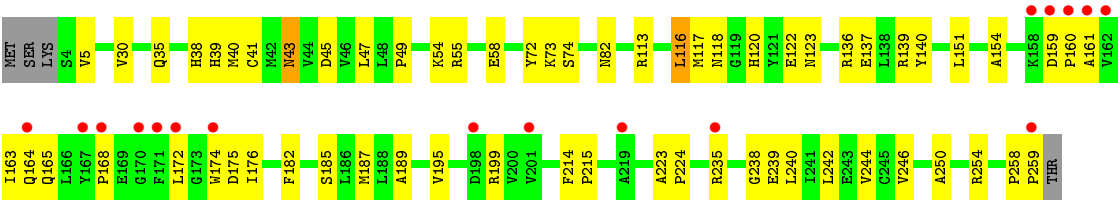


- Molecule 1: Creatinine amidohydrolase

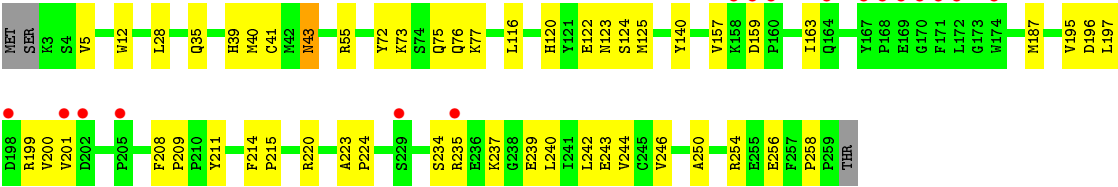
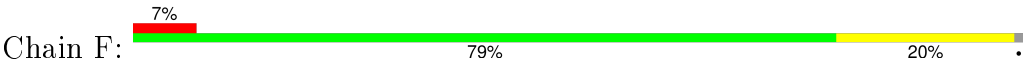


- Molecule 1: Creatinine amidohydrolase





• Molecule 1: Creatinine amidohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.80Å 60.24Å 146.20Å 90.00° 100.30° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 39.60 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.0 (20.00-2.00) 97.9 (39.60-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.199 , 0.231 0.200 , 0.231	Depositor DCC
R_{free} test set	6109 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	22.5	Xtriage
Anisotropy	0.595	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 121679 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12628	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% 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, 0.375 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: ZN, MN, 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	0.33	0/2026	0.59	0/2756
1	B	0.32	0/2017	0.59	0/2745
1	C	0.31	0/2013	0.58	0/2741
1	D	0.32	0/2022	0.58	0/2752
1	E	0.29	0/2006	0.55	0/2731
1	F	0.29	0/2026	0.54	0/2756
All	All	0.31	0/12110	0.57	0/16481

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	1979	0	1964	16	0
1	B	1970	0	1951	15	0
1	C	1966	0	1940	26	0
1	D	1975	0	1953	26	0
1	E	1960	0	1940	45	0
1	F	1979	0	1964	33	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	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
5	A	157	0	0	0	0
5	B	149	0	0	1	0
5	C	127	0	0	1	0
5	D	153	0	0	1	0
5	E	86	0	0	3	0
5	F	109	0	0	1	0
All	All	12628	0	11712	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:F:28:LEU:HD23	1:F:116:LEU:HD21	1.62	0.80
1:E:120:HIS:HB3	1:E:123:ASN:ND2	2.00	0.77
1:F:41:CYS:SG	1:F:43:ASN:ND2	2.65	0.69
1:F:120:HIS:HB3	1:F:123:ASN:ND2	2.08	0.68
1:C:74:SER:H	1:C:82:ASN:HD22	1.42	0.68
1:B:235:ARG:O	1:B:239:GLU:HG2	1.95	0.67
1:E:136:ARG:HA	1:E:139:ARG:NH1	2.10	0.66
1:A:3:LYS:HG3	1:A:4:SER:H	1.60	0.66
1:E:41:CYS:SG	1:E:43:ASN:ND2	2.70	0.65
1:F:256:GLU:C	1:F:258:PRO:HD3	2.18	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:11:THR:OG1	1:D:14:GLU:HG3	1.98	0.63
1:A:41:CYS:SG	1:A:43:ASN:ND2	2.71	0.63
1:E:161:ALA:O	1:E:165:GLN:HG3	1.99	0.63
1:B:250:ALA:O	1:B:254:ARG:HG3	1.99	0.62
1:E:74:SER:H	1:E:82:ASN:HD22	1.45	0.62
1:F:250:ALA:O	1:F:254:ARG:HG3	1.99	0.62
1:C:41:CYS:SG	1:C:43:ASN:ND2	2.73	0.61
1:D:43:ASN:C	1:D:43:ASN:HD22	2.03	0.61
1:B:120:HIS:HB3	1:B:123:ASN:ND2	2.16	0.61
1:B:160:PRO:O	1:B:164:GLN:HG2	2.00	0.61
1:D:159:ASP:OD2	1:D:161:ALA:HB3	2.01	0.61
1:E:159:ASP:OD2	1:E:161:ALA:HB3	2.02	0.60
1:D:41:CYS:SG	1:D:43:ASN:ND2	2.74	0.60
1:F:223:ALA:HB3	1:F:224:PRO:HD3	1.83	0.60
1:D:120:HIS:HB3	1:D:123:ASN:ND2	2.18	0.58
1:C:240:LEU:O	1:C:244:VAL:HG23	2.04	0.58
1:C:120:HIS:HB3	1:C:123:ASN:ND2	2.19	0.57
1:F:77:LYS:HE2	1:F:208:PHE:HB2	1.87	0.57
1:A:235:ARG:O	1:A:239:GLU:HG3	2.04	0.56
1:C:196:ASP:OD2	1:C:199:ARG:HG3	2.05	0.56
1:F:76:GLN:NE2	1:F:220:ARG:HB2	2.21	0.56
1:D:74:SER:H	1:D:82:ASN:HD22	1.52	0.56
1:B:41:CYS:SG	1:B:43:ASN:ND2	2.78	0.56
1:A:137:GLU:HG2	1:E:140:TYR:OH	2.06	0.55
1:E:159:ASP:O	1:E:163:ILE:HG12	2.06	0.55
1:E:240:LEU:O	1:E:244:VAL:HG23	2.06	0.55
1:E:189:ALA:HB1	1:E:239:GLU:OE2	2.06	0.55
1:E:250:ALA:O	1:E:254:ARG:HG3	2.06	0.54
1:E:154:ALA:HB1	1:E:174:TRP:CZ3	2.42	0.54
1:E:136:ARG:HA	1:E:139:ARG:HH11	1.70	0.54
1:C:38:HIS:CE1	1:C:224:PRO:HA	2.41	0.54
1:A:120:HIS:HB3	1:A:123:ASN:ND2	2.23	0.54
1:A:187:MET:HB3	1:A:195:VAL:HG21	1.90	0.54
1:E:214:PHE:HA	1:E:215:PRO:C	2.29	0.52
1:C:116:LEU:HD13	1:C:131:ILE:HD11	1.92	0.52
1:C:118:ASN:ND2	1:C:120:HIS:H	2.06	0.52
1:F:242:LEU:O	1:F:246:VAL:HG23	2.10	0.51
1:D:187:MET:HB3	1:D:195:VAL:HG21	1.91	0.51
1:A:159:ASP:HB3	1:A:162:VAL:HG23	1.92	0.51
1:C:188:LEU:O	1:C:192:PRO:HG3	2.11	0.51
1:E:118:ASN:HD21	1:E:123:ASN:HD22	1.58	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:240:LEU:O	1:F:244:VAL:HG23	2.10	0.51
1:D:114:LEU:HG	1:D:116:LEU:HD12	1.93	0.51
1:E:30:VAL:HG21	1:E:116:LEU:HD23	1.94	0.49
1:D:113:ARG:HD2	5:D:1606:HOH:O	2.12	0.49
1:E:164:GLN:NE2	1:E:168:PRO:HA	2.27	0.49
1:A:12:TRP:CD1	1:F:35:GLN:HB2	2.48	0.49
1:E:54:LYS:NZ	1:E:58:GLU:OE1	2.45	0.49
1:E:38:HIS:CE1	1:E:224:PRO:HA	2.48	0.48
1:A:214:PHE:HA	1:A:215:PRO:C	2.34	0.48
1:D:214:PHE:HA	1:D:215:PRO:C	2.34	0.48
1:F:5:VAL:HG22	5:F:1706:HOH:O	2.12	0.48
1:D:137:GLU:HG2	1:F:140:TYR:OH	2.12	0.48
1:D:159:ASP:HB3	1:D:162:VAL:HG23	1.95	0.48
1:A:187:MET:HB3	1:A:195:VAL:CG2	2.44	0.48
1:D:235:ARG:O	1:D:239:GLU:HG3	2.14	0.48
1:E:74:SER:H	1:E:82:ASN:ND2	2.12	0.48
1:F:214:PHE:HA	1:F:215:PRO:C	2.34	0.48
1:B:214:PHE:HA	1:B:215:PRO:C	2.34	0.47
1:F:234:SER:OG	1:F:237:LYS:HG3	2.14	0.47
1:C:26:LEU:HD23	1:C:63:LEU:HB2	1.96	0.47
1:F:55:ARG:HH11	1:F:55:ARG:HG3	1.79	0.46
1:D:39:HIS:CD2	1:D:40:MET:HG3	2.51	0.46
1:F:196:ASP:OD2	1:F:199:ARG:HG3	2.15	0.46
5:E:1171:HOH:O	1:F:125:MET:HG3	2.15	0.46
1:C:43:ASN:HD22	1:C:43:ASN:C	2.18	0.46
1:E:43:ASN:HD22	1:E:43:ASN:C	2.17	0.46
1:E:242:LEU:O	1:E:246:VAL:HG23	2.14	0.46
1:E:172:LEU:N	1:E:172:LEU:HD12	2.30	0.46
1:C:214:PHE:HA	1:C:215:PRO:C	2.36	0.46
1:F:39:HIS:CD2	1:F:40:MET:HG3	2.51	0.46
1:D:187:MET:HB3	1:D:195:VAL:CG2	2.46	0.46
1:B:251:ASP:OD1	1:B:254:ARG:NH1	2.49	0.46
1:D:43:ASN:C	1:D:43:ASN:ND2	2.68	0.46
1:E:55:ARG:HG3	1:E:55:ARG:HH11	1.81	0.46
1:E:43:ASN:HD21	1:E:47:LEU:HD11	1.81	0.45
1:E:117:MET:HE2	1:E:151:LEU:HD11	1.99	0.45
1:C:197:LEU:HA	1:C:200:VAL:HG23	1.99	0.45
1:F:159:ASP:O	1:F:163:ILE:HG12	2.17	0.45
1:E:223:ALA:N	1:E:224:PRO:HD2	2.32	0.45
1:C:55:ARG:HH11	1:C:55:ARG:HG3	1.82	0.45
1:A:125:MET:HG3	5:B:1038:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:118:ASN:ND2	1:D:120:HIS:H	2.14	0.44
1:C:258:PRO:HA	1:C:259:PRO:HD3	1.90	0.44
1:E:154:ALA:HB1	1:E:174:TRP:CH2	2.53	0.44
1:E:174:TRP:C	1:E:176:ILE:H	2.21	0.44
1:D:118:ASN:HD21	1:D:123:ASN:HD22	1.64	0.44
1:F:75:GLN:HG2	1:F:122:GLU:HG2	1.99	0.44
1:E:118:ASN:ND2	1:E:120:HIS:H	2.15	0.44
1:D:43:ASN:HD21	1:D:47:LEU:CD1	2.31	0.44
1:F:122:GLU:OE1	1:F:122:GLU:N	2.48	0.44
1:C:153:TYR:HE1	1:C:182:PHE:HE1	1.66	0.43
1:B:43:ASN:C	1:B:43:ASN:HD22	2.20	0.43
1:E:122:GLU:N	1:E:122:GLU:OE1	2.48	0.43
1:F:235:ARG:HG2	1:F:239:GLU:OE1	2.17	0.43
1:E:72:TYR:CG	1:E:73:LYS:N	2.86	0.43
1:A:35:GLN:HB2	1:F:12:TRP:CD1	2.54	0.43
1:D:12:TRP:CD1	1:E:35:GLN:HB2	2.54	0.43
1:E:235:ARG:HB3	1:E:235:ARG:NH2	2.34	0.42
1:E:258:PRO:HA	1:E:259:PRO:HD3	1.90	0.42
1:E:39:HIS:CD2	1:E:40:MET:HG3	2.54	0.42
1:F:72:TYR:CG	1:F:73:LYS:N	2.87	0.42
1:F:187:MET:HB3	1:F:195:VAL:HG21	2.01	0.42
1:B:12:TRP:CE3	1:C:85:PRO:HD2	2.54	0.42
1:C:172:LEU:HD13	1:C:205:PRO:HG2	2.02	0.42
1:C:235:ARG:HH22	1:C:239:GLU:HG3	1.84	0.42
1:F:197:LEU:HA	1:F:200:VAL:HG23	2.02	0.42
1:E:113:ARG:NE	5:E:1411:HOH:O	2.53	0.42
1:B:157:VAL:HG12	1:B:157:VAL:O	2.19	0.42
1:B:258:PRO:HA	1:B:259:PRO:HD3	1.88	0.42
1:D:43:ASN:HD21	1:D:47:LEU:HD11	1.85	0.42
1:A:159:ASP:HB3	1:A:162:VAL:CG2	2.49	0.41
1:D:37:GLY:HA2	1:D:224:PRO:O	2.20	0.41
1:F:201:VAL:HG11	1:F:224:PRO:HG2	2.02	0.41
1:A:243:GLU:HA	1:A:243:GLU:OE2	2.20	0.41
1:C:15:TYR:OH	1:C:109:HIS:HD2	2.02	0.41
1:E:5:VAL:HG22	5:E:1465:HOH:O	2.21	0.41
1:F:55:ARG:NH1	1:F:55:ARG:HG3	2.36	0.41
1:E:45:ASP:O	1:E:49:PRO:HG3	2.21	0.41
1:B:87:THR:O	1:C:109:HIS:HE1	2.03	0.41
1:E:199:ARG:HH21	1:E:199:ARG:HG2	1.85	0.41
5:C:1086:HOH:O	1:D:125:MET:HG3	2.20	0.41
1:B:141:ALA:HB2	1:D:140:TYR:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:223:ALA:HB3	1:C:224:PRO:HD3	2.02	0.41
1:C:203:HIS:HB2	1:C:204:PRO:HD2	2.02	0.41
1:C:74:SER:H	1:C:82:ASN:ND2	2.13	0.41
1:A:247:GLN:NE2	1:A:251:ASP:OD2	2.51	0.41
1:F:43:ASN:C	1:F:43:ASN:HD22	2.24	0.41
1:F:157:VAL:O	1:F:163:ILE:HD11	2.21	0.41
1:D:242:LEU:O	1:D:246:VAL:HG23	2.20	0.41
1:F:209:PRO:HB2	1:F:211:TYR:CE2	2.56	0.41
1:E:187:MET:HB3	1:E:195:VAL:HG21	2.02	0.41
1:A:75:GLN:HG2	1:A:122:GLU:HG2	2.03	0.41
1:B:159:ASP:HB3	1:B:162:VAL:HG23	2.02	0.41
1:C:43:ASN:HD21	1:C:47:LEU:HD11	1.86	0.40
1:F:76:GLN:HE22	1:F:220:ARG:HB2	1.85	0.40
1:C:140:TYR:OH	1:E:137:GLU:HG2	2.21	0.40
1:E:159:ASP:HA	1:E:160:PRO:HD3	1.99	0.40
1:E:174:TRP:HH2	1:E:182:PHE:CE2	2.40	0.40
1:B:114:LEU:HG	1:B:116:LEU:HD12	2.02	0.40
1:E:185:SER:O	1:E:238:GLY:HA3	2.21	0.40
1:D:114:LEU:CD2	1:D:116:LEU:HD11	2.51	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	255/260 (98%)	248 (97%)	7 (3%)	0	100	100
1	B	254/260 (98%)	249 (98%)	5 (2%)	0	100	100
1	C	254/260 (98%)	246 (97%)	8 (3%)	0	100	100
1	D	255/260 (98%)	248 (97%)	6 (2%)	1 (0%)	39	33
1	E	254/260 (98%)	247 (97%)	6 (2%)	1 (0%)	39	33

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	255/260 (98%)	250 (98%)	4 (2%)	1 (0%)	39	33
All	All	1527/1560 (98%)	1488 (97%)	36 (2%)	3 (0%)	52	48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	175	ASP
1	D	124	SER
1	F	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	211/214 (99%)	210 (100%)	1 (0%)	92	94
1	B	210/214 (98%)	207 (99%)	3 (1%)	74	77
1	C	209/214 (98%)	205 (98%)	4 (2%)	65	67
1	D	210/214 (98%)	208 (99%)	2 (1%)	82	85
1	E	208/214 (97%)	206 (99%)	2 (1%)	82	85
1	F	211/214 (99%)	209 (99%)	2 (1%)	84	88
All	All	1259/1284 (98%)	1245 (99%)	14 (1%)	80	83

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	B	43	ASN
1	B	212	ASP
1	B	247	GLN
1	C	38	HIS
1	C	43	ASN
1	C	116	LEU
1	C	212	ASP

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Mol	Chain	Res	Type
1	D	38	HIS
1	D	43	ASN
1	E	43	ASN
1	E	116	LEU
1	F	43	ASN
1	F	243	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	75	GLN
1	A	118	ASN
1	B	43	ASN
1	C	38	HIS
1	C	43	ASN
1	C	82	ASN
1	C	109	HIS
1	C	118	ASN
1	C	144	GLN
1	C	165	GLN
1	D	38	HIS
1	D	43	ASN
1	D	69	GLN
1	D	75	GLN
1	D	82	ASN
1	D	118	ASN
1	D	164	GLN
1	E	38	HIS
1	E	43	ASN
1	E	75	GLN
1	E	82	ASN
1	E	118	ASN
1	E	144	GLN
1	E	164	GLN
1	E	165	GLN
1	F	43	ASN
1	F	118	ASN

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, 18 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 [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	257/260 (98%)	-0.12	3 (1%) 81 81	10, 21, 36, 51	0
1	B	256/260 (98%)	-0.17	6 (2%) 64 64	10, 22, 39, 52	0
1	C	256/260 (98%)	-0.12	4 (1%) 74 75	11, 25, 47, 60	0
1	D	257/260 (98%)	-0.16	5 (1%) 70 70	13, 23, 40, 54	0
1	E	256/260 (98%)	0.34	17 (6%) 22 22	14, 32, 61, 71	0
1	F	257/260 (98%)	0.24	17 (6%) 22 22	14, 32, 64, 75	0
All	All	1539/1560 (98%)	0.00	52 (3%) 49 50	10, 25, 48, 75	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	161	ALA	5.2
1	E	170	GLY	4.7
1	C	171	PHE	4.4
1	E	174	TRP	4.3
1	F	202	ASP	4.0
1	F	164	GLN	3.9
1	F	174	TRP	3.8
1	E	158	LYS	3.7
1	E	160	PRO	3.7
1	D	169	GLU	3.6
1	E	164	GLN	3.5
1	F	158	LYS	3.5
1	B	172	LEU	3.4
1	F	168	PRO	3.4
1	B	169	GLU	3.3
1	A	171	PHE	3.3
1	E	171	PHE	3.2
1	E	259	PRO	3.1
1	F	171	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	F	170	GLY	3.1
1	E	168	PRO	3.0
1	B	164	GLN	3.0
1	E	201	VAL	3.0
1	B	171	PHE	3.0
1	E	235	ARG	2.8
1	F	205	PRO	2.7
1	C	168	PRO	2.7
1	E	172	LEU	2.7
1	F	159	ASP	2.6
1	F	169	GLU	2.6
1	E	167	TYR	2.5
1	A	169	GLU	2.5
1	F	167	TYR	2.5
1	D	171	PHE	2.5
1	F	229	SER	2.5
1	D	168	PRO	2.5
1	F	235	ARG	2.3
1	B	160	PRO	2.3
1	F	160	PRO	2.3
1	F	172	LEU	2.3
1	E	219	ALA	2.2
1	A	168	PRO	2.2
1	D	172	LEU	2.2
1	C	164	GLN	2.1
1	E	159	ASP	2.1
1	F	198	ASP	2.1
1	B	170	GLY	2.1
1	F	201	VAL	2.1
1	E	198	ASP	2.1
1	D	170	GLY	2.1
1	C	202	ASP	2.0
1	E	162	VAL	2.0

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

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

6.3 Carbohydrates ⓘ

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	CL	A	302	1/1	0.95	0.23	4.60	33,33,33,33	0
4	CL	F	307	1/1	0.89	0.18	2.90	47,47,47,47	0
4	CL	D	305	1/1	0.96	0.17	2.82	35,35,35,35	0
4	CL	B	303	1/1	0.94	0.15	1.46	31,31,31,31	0
4	CL	C	304	1/1	0.98	0.14	0.78	29,29,29,29	0
2	MN	D	300	1/1	0.99	0.12	0.66	38,38,38,38	0
4	CL	E	306	1/1	0.94	0.16	0.63	37,37,37,37	0
2	MN	F	300	1/1	0.96	0.12	0.58	52,52,52,52	0
2	MN	E	300	1/1	0.94	0.15	0.55	54,54,54,54	0
2	MN	A	300	1/1	0.99	0.12	0.53	36,36,36,36	0
2	MN	B	300	1/1	0.99	0.10	-0.47	39,39,39,39	0
2	MN	C	300	1/1	0.99	0.09	-0.75	44,44,44,44	0
3	ZN	A	301	1/1	0.99	0.05	-2.30	33,33,33,33	0
3	ZN	E	301	1/1	0.97	0.07	-2.66	45,45,45,45	0
3	ZN	B	301	1/1	0.98	0.04	-2.69	33,33,33,33	0
3	ZN	C	301	1/1	0.97	0.06	-2.82	37,37,37,37	0
3	ZN	F	301	1/1	0.95	0.04	-3.78	45,45,45,45	0
3	ZN	D	301	1/1	0.98	0.05	-5.44	35,35,35,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.