



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 12:47 PM GMT

PDB ID : 3A6G
Title : W154F 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 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/ValidationPDFNotes.html>

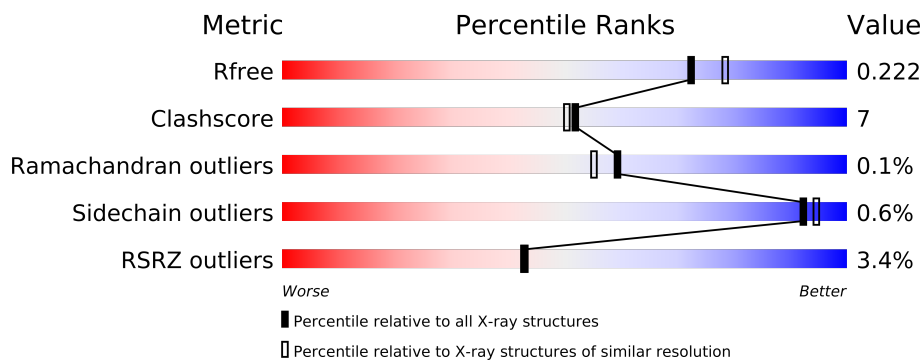
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	260	
1	B	260	
1	C	260	
1	D	260	
1	E	260	
1	F	260	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12321 atoms, of which 0 are hydrogen and 0 are deuterium.

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
			1940	1238	332	359	11			
1	B	253	Total	C	N	O	S	0	0	0
			1926	1229	328	358	11			
1	C	254	Total	C	N	O	S	0	0	0
			1944	1244	330	359	11			
1	D	257	Total	C	N	O	S	0	0	0
			1981	1269	336	365	11			
1	E	257	Total	C	N	O	S	0	0	0
			1981	1269	336	365	11			
1	F	257	Total	C	N	O	S	0	0	0
			1977	1266	335	365	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	154	PHE	TRP	ENGINEERED	UNP P83772
B	154	PHE	TRP	ENGINEERED	UNP P83772
C	154	PHE	TRP	ENGINEERED	UNP P83772
D	154	PHE	TRP	ENGINEERED	UNP P83772
E	154	PHE	TRP	ENGINEERED	UNP P83772
F	154	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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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).

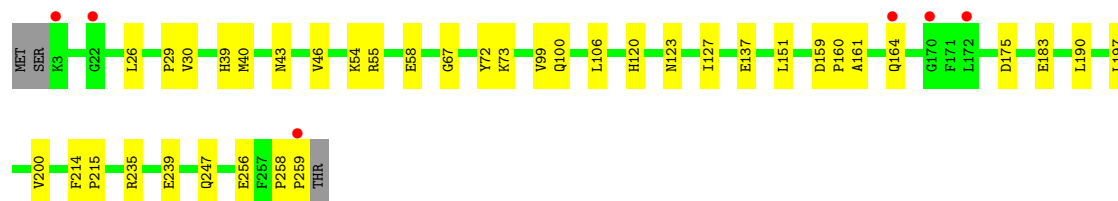
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total 1	Zn 1	0	0
3	E	1	Total 1	Zn 1	0	0
3	B	1	Total 1	Zn 1	0	0
3	C	1	Total 1	Zn 1	0	0
3	A	1	Total 1	Zn 1	0	0
3	F	1	Total 1	Zn 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	97	Total 97	O 97	0	0
4	B	96	Total 96	O 96	0	0
4	C	93	Total 93	O 93	0	0
4	D	98	Total 98	O 98	0	0
4	E	95	Total 95	O 95	0	0
4	F	81	Total 81	O 81	0	0

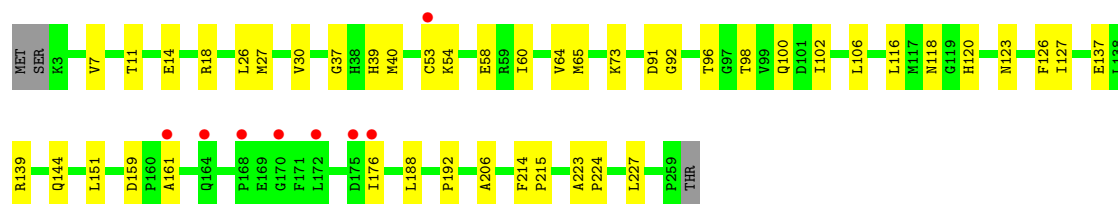
- Molecule 1: Creatinine amidohydrolase

Chain E:



- Molecule 1: Creatinine amidohydrolase

Chain F:



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	164.50Å 164.50Å 164.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.00 45.62 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.00) 100.0 (45.62-2.00)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.199 , 0.216 0.203 , 0.222	Depositor DCC
R_{free} test set	8621 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.319	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 46.0	EDS
Estimated twinning fraction	0.018 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 173055 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12321	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.31	0/1985	0.57	0/2702
1	B	0.31	0/1970	0.57	0/2681
1	C	0.31	0/1989	0.56	0/2705
1	D	0.31	0/2029	0.57	0/2760
1	E	0.31	0/2029	0.57	0/2760
1	F	0.32	0/2025	0.56	0/2756
All	All	0.31	0/12027	0.57	0/16364

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1940	0	1897	33	0
1	B	1926	0	1891	22	0
1	C	1944	0	1920	25	0
1	D	1981	0	1964	27	0
1	E	1981	0	1964	25	0
1	F	1977	0	1953	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	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	97	0	0	2	0
4	B	96	0	0	1	0
4	C	93	0	0	3	0
4	D	98	0	0	1	0
4	E	95	0	0	2	0
4	F	81	0	0	1	0
All	All	12321	0	11589	168	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (168) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:35:GLN:H	1:A:87:THR:HG22	1.33	0.92
1:D:60:ILE:HD11	1:D:250:ALA:HB1	1.53	0.90
1:F:176:ILE:HD11	1:F:206:ALA:HB3	1.64	0.78
1:A:87:THR:HG21	4:A:1066:HOH:O	1.84	0.75
1:F:53:CYS:SG	1:F:64:VAL:HG22	2.28	0.74
1:C:54:LYS:O	1:C:58:GLU:HG3	1.89	0.72
1:E:55:ARG:HG2	1:E:55:ARG:HH11	1.54	0.71
1:E:197:LEU:O	1:E:200:VAL:HG12	1.91	0.70
1:A:176:ILE:HD13	1:A:205:PRO:HB3	1.74	0.69
1:D:54:LYS:O	1:D:58:GLU:HG3	1.93	0.69
1:F:7:VAL:HG21	1:F:65:MET:HE2	1.73	0.68
1:E:30:VAL:HG11	1:E:127:ILE:HD11	1.76	0.68
1:F:188:LEU:O	1:F:192:PRO:HG3	1.94	0.67
1:A:176:ILE:HD11	1:A:206:ALA:N	2.11	0.66
1:C:235:ARG:O	1:C:239:GLU:HG2	1.94	0.66
1:D:72:TYR:H	1:D:123:ASN:ND2	1.94	0.65
1:A:235:ARG:O	1:A:239:GLU:HG3	1.97	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:35:GLN:H	1:A:87:THR:CG2	2.09	0.63
1:B:28:LEU:HD23	1:B:116:LEU:HD21	1.81	0.63
1:A:176:ILE:HD11	1:A:206:ALA:H	1.66	0.60
1:F:159:ASP:OD2	1:F:161:ALA:HB3	2.02	0.60
1:F:26:LEU:HD13	1:F:65:MET:HE3	1.85	0.59
1:F:27:MET:CE	1:F:60:ILE:HD11	2.33	0.59
1:D:26:LEU:HD23	1:D:106:LEU:HD22	1.84	0.59
1:A:35:GLN:N	1:A:87:THR:HG22	2.12	0.58
1:D:249:ILE:O	1:D:253:ILE:HG12	2.04	0.57
1:B:198:ASP:HB2	4:B:1486:HOH:O	2.04	0.57
1:F:27:MET:HB2	1:F:53:CYS:SG	2.44	0.57
1:B:159:ASP:OD2	1:B:161:ALA:HB3	2.04	0.57
1:F:54:LYS:O	1:F:58:GLU:HG3	2.05	0.57
1:A:54:LYS:O	1:A:58:GLU:HG3	2.04	0.57
1:B:235:ARG:O	1:B:239:GLU:HG3	2.04	0.57
1:D:26:LEU:CD2	1:D:106:LEU:HD22	2.34	0.56
1:C:120:HIS:HB3	1:C:123:ASN:ND2	2.21	0.56
1:F:139:ARG:HD3	1:F:144:GLN:HE21	1.71	0.56
1:A:28:LEU:HD23	1:A:116:LEU:CD2	2.36	0.56
1:D:19:VAL:HG13	4:D:1406:HOH:O	2.05	0.56
1:B:249:ILE:O	1:B:253:ILE:HG12	2.06	0.55
1:F:120:HIS:HB3	1:F:123:ASN:ND2	2.22	0.55
1:A:56:VAL:O	1:A:60:ILE:HG12	2.07	0.55
1:E:55:ARG:NH1	1:E:55:ARG:HG2	2.22	0.55
1:F:30:VAL:CG1	1:F:118:ASN:HA	2.37	0.55
1:E:54:LYS:O	1:E:58:GLU:HG3	2.08	0.54
1:E:99:VAL:HG21	1:E:127:ILE:HD13	1.89	0.54
1:D:28:LEU:HD11	1:D:68:LEU:CD1	2.38	0.54
1:F:11:THR:OG1	1:F:14:GLU:HG3	2.08	0.54
1:D:62:ALA:C	1:D:63:LEU:HD12	2.28	0.54
1:B:176:ILE:HD11	1:B:206:ALA:N	2.22	0.54
1:A:98:THR:O	1:A:102:ILE:HG12	2.08	0.53
1:A:176:ILE:HG13	1:A:227:LEU:HB3	1.91	0.53
1:D:214:PHE:HA	1:D:215:PRO:C	2.29	0.52
1:E:55:ARG:HH21	1:E:190:LEU:CD2	2.23	0.52
1:B:54:LYS:O	1:B:58:GLU:HG3	2.10	0.52
1:C:155:ASP:O	1:C:158:LYS:HD2	2.09	0.52
1:E:120:HIS:HB3	1:E:123:ASN:ND2	2.25	0.52
1:D:39:HIS:CD2	1:D:40:MET:HG3	2.45	0.52
1:F:30:VAL:HG13	1:F:118:ASN:HA	1.91	0.51
1:F:27:MET:HE3	1:F:60:ILE:HD11	1.92	0.51
1:E:256:GLU:C	1:E:258:PRO:HD3	2.31	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:214:PHE:HA	1:B:215:PRO:C	2.32	0.51
1:A:28:LEU:HD23	1:A:116:LEU:HD21	1.93	0.50
1:F:30:VAL:O	1:F:30:VAL:HG13	2.12	0.50
1:E:43:ASN:O	1:E:46:VAL:HG12	2.12	0.50
1:F:18:ARG:HD3	4:F:1352:HOH:O	2.11	0.50
1:D:35:GLN:NE2	1:D:37:GLY:H	2.10	0.49
1:B:235:ARG:HG2	1:B:235:ARG:HH21	1.76	0.49
1:F:30:VAL:HG11	1:F:127:ILE:HD11	1.94	0.49
1:C:23:ASP:HB2	1:C:113:ARG:NH2	2.26	0.49
1:F:39:HIS:CD2	1:F:40:MET:HG3	2.47	0.49
1:A:120:HIS:HB3	1:A:123:ASN:ND2	2.28	0.49
1:B:28:LEU:HD23	1:B:116:LEU:CD2	2.42	0.49
1:B:176:ILE:HD11	1:B:206:ALA:H	1.77	0.49
1:B:11:THR:OG1	1:B:14:GLU:HG3	2.12	0.49
1:D:24:CYS:SG	1:D:63:LEU:HD13	2.53	0.49
1:C:55:ARG:HH11	1:C:55:ARG:HG3	1.78	0.49
1:F:96:THR:HG22	1:F:126:PHE:O	2.13	0.49
1:E:214:PHE:HA	1:E:215:PRO:C	2.33	0.49
1:B:120:HIS:HB3	1:B:123:ASN:ND2	2.28	0.48
1:C:214:PHE:HA	1:C:215:PRO:C	2.33	0.48
1:A:39:HIS:CD2	1:A:40:MET:HG3	2.48	0.48
1:F:73:LYS:NZ	1:F:91:ASP:HA	2.28	0.48
1:A:34:GLU:HA	1:A:87:THR:HG22	1.96	0.48
1:E:159:ASP:OD2	1:E:161:ALA:HB3	2.13	0.48
1:C:153:TYR:HE1	1:C:182:PHE:HE1	1.61	0.47
1:B:56:VAL:O	1:B:60:ILE:HG12	2.14	0.47
1:F:151:LEU:HD12	1:F:151:LEU:C	2.35	0.47
1:F:214:PHE:HA	1:F:215:PRO:C	2.35	0.47
1:F:116:LEU:HD22	1:F:127:ILE:HG23	1.96	0.47
1:C:212:ASP:HB2	1:D:150:VAL:HG13	1.95	0.47
1:A:214:PHE:HA	1:A:215:PRO:C	2.35	0.47
1:A:75:GLN:HG2	1:A:122:GLU:HG2	1.96	0.47
1:F:37:GLY:HA2	1:F:224:PRO:O	2.15	0.46
1:E:183:GLU:HG3	4:E:1186:HOH:O	2.16	0.46
1:C:37:GLY:HA2	1:C:224:PRO:O	2.16	0.46
1:C:256:GLU:C	1:C:258:PRO:HD3	2.36	0.46
1:E:100:GLN:NE2	1:E:137:GLU:OE2	2.48	0.46
1:B:151:LEU:C	1:B:151:LEU:HD12	2.36	0.46
1:A:153:TYR:HE1	1:A:182:PHE:HE1	1.63	0.46
1:C:39:HIS:CD2	1:C:40:MET:HG3	2.50	0.46
1:C:152:SER:HB2	1:C:155:ASP:OD2	2.16	0.46
1:F:223:ALA:HB3	1:F:224:PRO:HD3	1.98	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:11:THR:OG1	1:A:14:GLU:HG3	2.17	0.45
1:D:55:ARG:HH11	1:D:55:ARG:HG3	1.81	0.45
1:B:139:ARG:HG3	1:B:144:GLN:HE21	1.81	0.45
1:C:242:LEU:O	1:C:246:VAL:HG23	2.17	0.45
1:E:160:PRO:O	1:E:164:GLN:HG3	2.16	0.45
1:F:27:MET:CB	1:F:53:CYS:SG	3.05	0.45
1:A:83:HIS:H	1:A:83:HIS:CD2	2.34	0.45
1:F:92:GLY:O	1:F:96:THR:HG23	2.17	0.45
1:C:40:MET:HG2	1:C:195:VAL:HG22	1.98	0.45
1:D:11:THR:OG1	1:D:14:GLU:HG3	2.17	0.44
1:D:159:ASP:OD2	1:D:162:VAL:HG23	2.17	0.44
1:E:247:GLN:NE2	4:E:1510:HOH:O	2.51	0.44
1:E:39:HIS:CD2	1:E:40:MET:HG3	2.52	0.44
1:C:111:ALA:HB1	4:C:1502:HOH:O	2.17	0.44
1:D:242:LEU:O	1:D:246:VAL:HG23	2.17	0.44
1:A:241:ILE:O	1:A:244:VAL:HG12	2.17	0.43
1:E:235:ARG:NH1	1:E:239:GLU:OE2	2.51	0.43
1:C:183:GLU:HG3	4:C:1473:HOH:O	2.16	0.43
1:C:139:ARG:HD3	1:C:144:GLN:HE21	1.82	0.43
1:D:258:PRO:HA	1:D:259:PRO:HD3	1.85	0.43
1:B:39:HIS:CD2	1:B:40:MET:HG3	2.53	0.43
1:B:176:ILE:HD13	1:B:205:PRO:HB3	2.01	0.43
1:D:75:GLN:HG2	1:D:122:GLU:HG2	2.01	0.43
1:A:100:GLN:NE2	1:A:137:GLU:OE2	2.50	0.43
1:A:77:LYS:HE2	1:A:208:PHE:HB3	2.00	0.43
1:B:223:ALA:HB3	1:B:224:PRO:HD3	2.01	0.43
1:F:98:THR:O	1:F:102:ILE:HG12	2.19	0.43
1:C:83:HIS:CD2	1:C:83:HIS:H	2.37	0.43
1:D:223:ALA:N	1:D:224:PRO:HD2	2.34	0.43
1:C:223:ALA:HB3	1:C:224:PRO:HD3	2.01	0.42
1:D:115:VAL:HG21	1:D:253:ILE:HD12	2.00	0.42
1:F:30:VAL:HG11	1:F:127:ILE:CD1	2.49	0.42
1:F:53:CYS:SG	1:F:64:VAL:CG2	3.03	0.42
1:F:53:CYS:SG	1:F:64:VAL:HG13	2.59	0.42
1:D:68:LEU:N	1:D:68:LEU:HD12	2.34	0.42
1:E:258:PRO:HA	1:E:259:PRO:HD3	1.92	0.42
1:A:83:HIS:HE1	4:A:1105:HOH:O	2.02	0.42
1:E:26:LEU:HD13	1:E:106:LEU:HD22	2.01	0.42
1:E:55:ARG:CG	1:E:55:ARG:NH1	2.83	0.42
1:F:100:GLN:NE2	1:F:137:GLU:OE2	2.50	0.42
1:E:72:TYR:CG	1:E:73:LYS:N	2.88	0.42
1:F:26:LEU:HD13	1:F:106:LEU:HD22	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:29:PRO:HG2	1:E:67:GLY:CA	2.48	0.42
1:C:159:ASP:OD2	1:C:162:VAL:HG23	2.20	0.42
1:C:83:HIS:HE1	4:C:1167:HOH:O	2.03	0.41
1:C:208:PHE:HB3	1:C:209:PRO:HD2	2.02	0.41
1:E:26:LEU:CD1	1:E:106:LEU:HD22	2.50	0.41
1:A:256:GLU:C	1:A:258:PRO:HD3	2.40	0.41
1:C:131:ILE:CD1	1:C:150:VAL:HG21	2.51	0.41
1:D:63:LEU:HD12	1:D:63:LEU:N	2.34	0.41
1:C:72:TYR:CG	1:C:73:LYS:N	2.89	0.41
1:A:26:LEU:HD13	1:A:106:LEU:HD22	2.03	0.41
1:B:176:ILE:HG13	1:B:227:LEU:HB3	2.02	0.41
1:D:166:LEU:C	1:D:168:PRO:HD3	2.40	0.41
1:F:176:ILE:CD1	1:F:227:LEU:HB3	2.50	0.41
1:A:176:ILE:CD1	1:A:205:PRO:HB3	2.48	0.41
1:B:239:GLU:O	1:B:243:GLU:HG3	2.21	0.41
1:A:26:LEU:CD1	1:A:106:LEU:HD22	2.51	0.41
1:A:72:TYR:CG	1:A:73:LYS:N	2.88	0.41
1:A:18:ARG:HD2	1:A:18:ARG:HA	1.88	0.41
1:F:176:ILE:HG13	1:F:227:LEU:O	2.21	0.40
1:A:250:ALA:O	1:A:254:ARG:HG3	2.21	0.40
1:D:71:GLY:HA3	1:D:123:ASN:HD21	1.86	0.40
1:E:151:LEU:C	1:E:151:LEU:HD12	2.42	0.40
1:D:72:TYR:H	1:D:123:ASN:HD21	1.64	0.40
1:B:176:ILE:HG12	1:B:176:ILE:O	2.21	0.40
1:F:73:LYS:HZ1	1:F:91:ASP:HA	1.86	0.40

There are no symmetry-related clashes.

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	255/260 (98%)	246 (96%)	9 (4%)	0	100	100
1	B	249/260 (96%)	240 (96%)	7 (3%)	2 (1%)	27	17
1	C	250/260 (96%)	243 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	255/260 (98%)	250 (98%)	5 (2%)	0	100	100
1	E	255/260 (98%)	249 (98%)	6 (2%)	0	100	100
1	F	255/260 (98%)	251 (98%)	4 (2%)	0	100	100
All	All	1519/1560 (97%)	1479 (97%)	38 (2%)	2 (0%)	59	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	176	ILE
1	B	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	201/215 (94%)	200 (100%)	1 (0%)	94	96
1	B	203/215 (94%)	202 (100%)	1 (0%)	94	96
1	C	206/215 (96%)	204 (99%)	2 (1%)	85	88
1	D	211/215 (98%)	208 (99%)	3 (1%)	78	81
1	E	211/215 (98%)	210 (100%)	1 (0%)	94	96
1	F	210/215 (98%)	210 (100%)	0	100	100
All	All	1242/1290 (96%)	1234 (99%)	8 (1%)	92	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	212	ASP
1	B	212	ASP
1	C	116	LEU
1	C	193	ASP
1	D	35	GLN
1	D	153	TYR
1	D	175	ASP
1	E	175	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	HIS
1	B	144	GLN
1	C	69	GLN
1	C	83	HIS
1	C	144	GLN
1	D	35	GLN
1	D	75	GLN
1	D	123	ASN
1	F	144	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 12 ligands modelled in this entry, 12 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.

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 ⓘ

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%)	-0.27	10 (3%) 37 37	10, 18, 44, 54	0
1	B	253/260 (97%)	-0.35	8 (3%) 45 45	10, 18, 46, 57	0
1	C	254/260 (97%)	-0.26	10 (3%) 37 37	10, 20, 46, 59	0
1	D	257/260 (98%)	-0.30	10 (3%) 37 37	11, 19, 43, 53	0
1	E	257/260 (98%)	-0.33	6 (2%) 57 57	11, 19, 39, 46	0
1	F	257/260 (98%)	-0.23	8 (3%) 47 46	10, 19, 46, 53	0
All	All	1535/1560 (98%)	-0.29	52 (3%) 43 43	10, 19, 44, 59	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	164	GLN	4.1
1	C	171	PHE	4.0
1	D	158	LYS	3.8
1	D	259	PRO	3.7
1	F	172	LEU	3.7
1	A	173	GLY	3.5
1	E	22	GLY	3.4
1	A	172	LEU	3.3
1	B	170	GLY	3.2
1	B	176	ILE	3.2
1	A	175	ASP	3.2
1	D	22	GLY	3.2
1	A	22	GLY	3.1
1	A	160	PRO	3.1
1	C	22	GLY	3.0
1	E	164	GLN	3.0
1	D	171	PHE	3.0
1	F	176	ILE	3.0
1	C	160	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	F	161	ALA	2.9
1	F	175	ASP	2.8
1	B	168	PRO	2.8
1	A	176	ILE	2.7
1	B	160	PRO	2.7
1	D	165	GLN	2.7
1	F	170	GLY	2.7
1	B	158	LYS	2.6
1	D	164	GLN	2.6
1	B	175	ASP	2.6
1	A	21	ALA	2.6
1	E	3	LYS	2.6
1	C	154	PHE	2.6
1	F	53	CYS	2.6
1	F	168	PRO	2.5
1	D	172	LEU	2.5
1	D	175	ASP	2.5
1	B	167	TYR	2.5
1	E	259	PRO	2.5
1	A	168	PRO	2.4
1	D	170	GLY	2.4
1	D	161	ALA	2.4
1	C	159	ASP	2.2
1	E	172	LEU	2.2
1	A	174	TRP	2.2
1	C	165	GLN	2.1
1	E	170	GLY	2.1
1	A	170	GLY	2.1
1	C	168	PRO	2.1
1	B	165	GLN	2.0
1	C	158	LYS	2.0
1	C	21	ALA	2.0
1	C	170	GLY	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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	ZN	A	301	1/1	0.05	-2.30	32,32,32,32	0
3	ZN	F	301	1/1	0.04	-2.48	30,30,30,30	0
2	MN	F	300	1/1	0.03	-2.85	26,26,26,26	0
3	ZN	D	301	1/1	0.04	-3.27	31,31,31,31	0
2	MN	C	300	1/1	0.03	-3.34	24,24,24,24	0
2	MN	B	300	1/1	0.03	-3.50	23,23,23,23	0
3	ZN	B	301	1/1	0.04	-3.56	30,30,30,30	0
2	MN	E	300	1/1	0.03	-3.79	22,22,22,22	0
2	MN	A	300	1/1	0.02	-4.64	23,23,23,23	0
3	ZN	C	301	1/1	0.02	-5.23	34,34,34,34	0
2	MN	D	300	1/1	0.02	-6.11	25,25,25,25	0
3	ZN	E	301	1/1	0.02	-6.81	30,30,30,30	0

6.5 Other polymers

There are no such residues in this entry.