



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 04:11 PM GMT

PDB ID : 3A6L  
Title : E122Q mutant creatininase, Zn-Zn type  
Authors : Nakajima, Y.; Yamashita, K.; Ito, K.; Yoshimoto, T.  
Deposited on : 2009-09-02  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

---

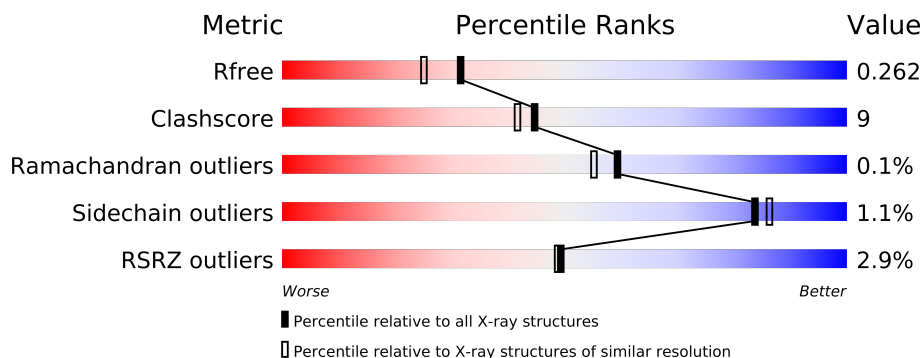
The following versions of software and data (see [references](#)) were used in the production of this report:

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  $\geq 3$ , 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	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	CL	A	302	-	X
3	CL	B	302	-	X

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Geometry	Electron density
3	CL	C	302	-	X
3	CL	D	302	-	X
3	CL	E	302	-	X
3	CL	F	302	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12469 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
			1976	1265	337	363	11			
1	B	257	Total	C	N	O	S	0	0	0
			1988	1273	338	366	11			
1	C	258	Total	C	N	O	S	0	0	0
			1990	1274	339	366	11			
1	D	256	Total	C	N	O	S	0	0	0
			1979	1267	336	365	11			
1	E	256	Total	C	N	O	S	0	0	0
			1975	1265	336	363	11			
1	F	257	Total	C	N	O	S	0	0	0
			1988	1273	338	366	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	122	GLN	GLU	ENGINEERED	UNP P83772
B	122	GLN	GLU	ENGINEERED	UNP P83772
C	122	GLN	GLU	ENGINEERED	UNP P83772
D	122	GLN	GLU	ENGINEERED	UNP P83772
E	122	GLN	GLU	ENGINEERED	UNP P83772
F	122	GLN	GLU	ENGINEERED	UNP P83772

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

*Continued on next page...*

*Continued from previous page...*

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

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 4 is water.

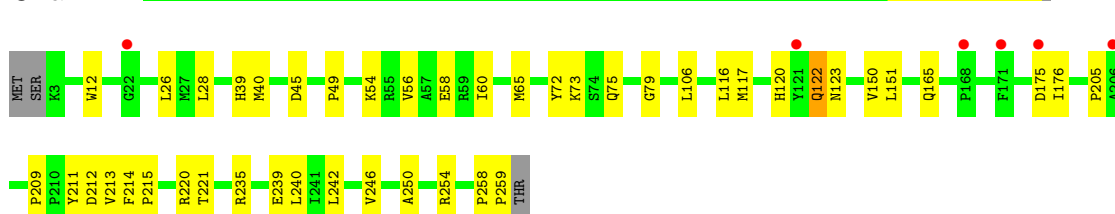
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	85	Total 85	O 85	0	0
4	B	88	Total 88	O 88	0	0
4	C	87	Total 87	O 87	0	0
4	D	103	Total 103	O 103	0	0
4	E	102	Total 102	O 102	0	0
4	F	90	Total 90	O 90	0	0

### 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.

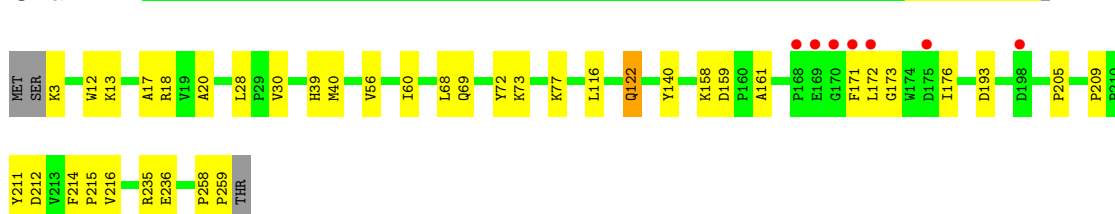
- Molecule 1: Creatinine amidohydrolase

Chain A:



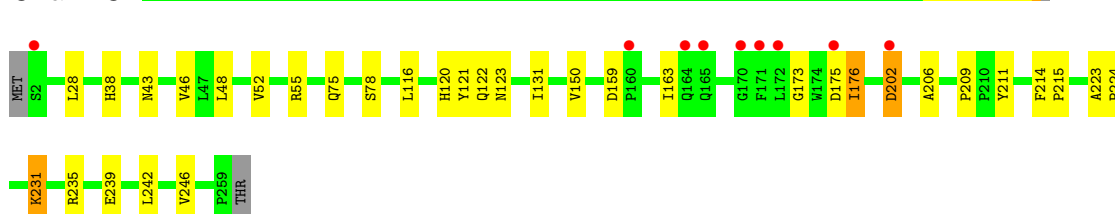
- Molecule 1: Creatinine amidohydrolase

Chain B:



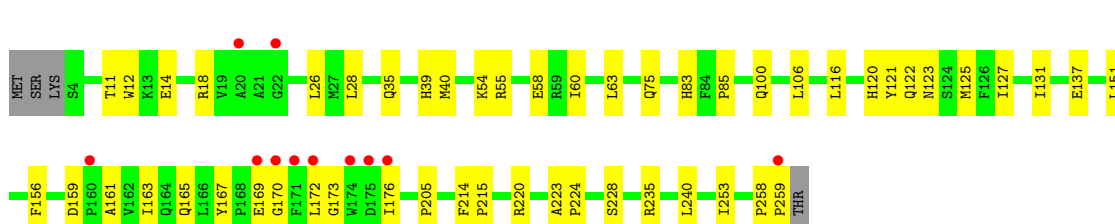
- Molecule 1: Creatinine amidohydrolase

Chain C:



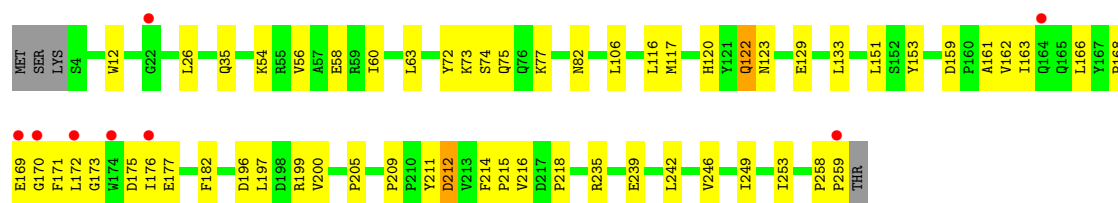
- Molecule 1: Creatinine amidohydrolase

Chain D:



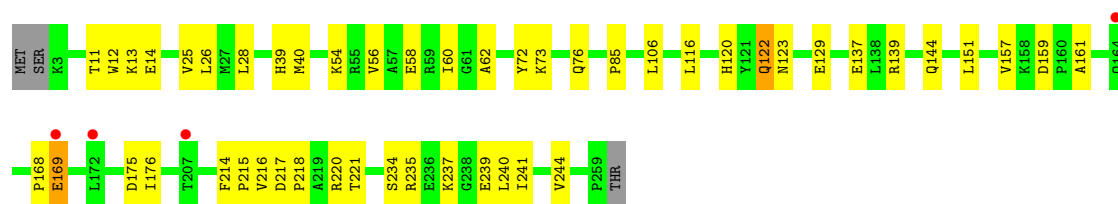
- Molecule 1: Creatinine amidohydrolase

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, $\alpha$ , $\beta$ , $\gamma$	164.40Å 164.40Å 164.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.00 32.89 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-2.00) 99.8 (32.89-2.00)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.05 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.235 , 0.259 0.237 , 0.262	Depositor DCC
$R_{free}$ test set	8630 reflections (5.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.6	Xtriage
Anisotropy	0.214	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 45.3	EDS
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 172480 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	12469	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.28% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities 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, 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	0.34	0/2025	0.59	0/2758
1	B	0.34	0/2037	0.59	0/2772
1	C	0.34	0/2039	0.58	0/2775
1	D	0.34	0/2028	0.59	0/2761
1	E	0.34	0/2024	0.59	0/2756
1	F	0.34	0/2037	0.58	0/2772
All	All	0.34	0/12190	0.59	0/16594

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	1976	0	1944	35	0
1	B	1988	0	1971	30	0
1	C	1990	0	1972	26	0
1	D	1979	0	1958	41	0
1	E	1975	0	1954	43	0
1	F	1988	0	1971	35	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	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	85	0	0	2	0
4	B	88	0	0	3	0
4	C	87	0	0	1	0
4	D	103	0	0	4	0
4	E	102	0	0	0	0
4	F	90	0	0	2	0
All	All	12469	0	11770	202	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 9.

All (202) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:122:GLN:H	1:A:122:GLN:NE2	1.49	1.09
1:A:122:GLN:N	1:A:122:GLN:HE21	1.59	0.99
1:C:75:GLN:HG2	1:C:122:GLN:HG2	1.49	0.91
1:D:75:GLN:H	1:D:122:GLN:HE21	1.18	0.89
1:A:75:GLN:HG2	1:A:122:GLN:HG2	1.52	0.89
1:D:176:ILE:O	1:D:228:SER:HB2	1.76	0.85
1:E:122:GLN:H	1:E:122:GLN:HE21	1.19	0.84
1:D:75:GLN:H	1:D:122:GLN:NE2	1.75	0.83
1:B:176:ILE:HD11	1:B:205:PRO:HB3	1.63	0.81
1:D:165:GLN:NE2	1:D:240:LEU:HD22	1.99	0.77
1:E:74:SER:H	1:E:82:ASN:HD22	1.29	0.77
1:D:75:GLN:HG2	1:D:122:GLN:HG2	1.68	0.75
1:B:235:ARG:HH11	1:B:236:GLU:HG3	1.53	0.73
1:E:75:GLN:HG2	1:E:122:GLN:HG2	1.69	0.72
1:C:176:ILE:H	1:C:176:ILE:HD13	1.56	0.71
1:C:120:HIS:HB3	1:C:123:ASN:ND2	2.06	0.71
1:D:176:ILE:HD12	1:D:205:PRO:HB3	1.73	0.70

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:122:GLN:H	1:E:122:GLN:NE2	1.89	0.69
1:C:43:ASN:O	1:C:46:VAL:HG22	1.92	0.68
1:D:120:HIS:HB3	1:D:123:ASN:ND2	2.08	0.68
1:E:176:ILE:HD11	1:E:205:PRO:HB3	1.76	0.68
1:E:159:ASP:OD2	1:E:161:ALA:HB3	1.93	0.67
1:B:173:GLY:O	1:B:176:ILE:HG22	1.94	0.67
1:B:235:ARG:NH1	1:B:236:GLU:HG3	2.09	0.67
1:E:120:HIS:HB3	1:E:123:ASN:ND2	2.09	0.67
1:A:122:GLN:H	1:A:122:GLN:HE21	0.76	0.65
1:D:26:LEU:HD13	1:D:106:LEU:HD22	1.79	0.65
1:B:13:LYS:HZ2	1:C:38:HIS:CE1	2.15	0.65
1:A:122:GLN:N	1:A:122:GLN:NE2	2.29	0.65
1:C:202:ASP:OD2	1:C:231:LYS:HG3	1.98	0.63
1:E:54:LYS:O	1:E:58:GLU:HG3	1.98	0.63
1:A:120:HIS:HB3	1:A:123:ASN:ND2	2.13	0.63
1:F:120:HIS:HB3	1:F:123:ASN:ND2	2.14	0.62
1:E:173:GLY:O	1:E:176:ILE:HG22	2.01	0.61
1:C:131:ILE:CD1	1:C:150:VAL:HG21	2.31	0.61
1:B:172:LEU:HB2	1:B:176:ILE:HG21	1.83	0.60
1:D:11:THR:OG1	1:D:14:GLU:HG3	2.01	0.60
1:D:28:LEU:HD23	1:D:116:LEU:HD21	1.82	0.60
1:D:220:ARG:HD3	4:D:1419:HOH:O	2.02	0.59
1:C:121:TYR:HD2	1:C:122:GLN:HE21	1.50	0.58
1:F:220:ARG:HD3	4:F:1272:HOH:O	2.02	0.58
1:F:120:HIS:NE2	1:F:122:GLN:HG2	2.18	0.58
1:E:74:SER:H	1:E:82:ASN:ND2	2.01	0.58
1:F:11:THR:OG1	1:F:14:GLU:HG3	2.03	0.58
1:D:169:GLU:HB2	4:D:1397:HOH:O	2.03	0.58
1:F:12:TRP:CZ2	1:F:13:LYS:HD3	2.39	0.57
1:C:235:ARG:O	1:C:239:GLU:HG2	2.05	0.57
1:C:159:ASP:O	1:C:163:ILE:HG12	2.05	0.57
1:D:75:GLN:N	1:D:122:GLN:HE21	1.96	0.57
1:F:25:VAL:HG12	4:F:1268:HOH:O	2.05	0.57
1:F:235:ARG:O	1:F:239:GLU:HG3	2.04	0.56
1:D:39:HIS:CD2	1:D:40:MET:HG3	2.40	0.56
1:F:54:LYS:O	1:F:58:GLU:HG3	2.05	0.56
1:F:240:LEU:O	1:F:244:VAL:HG23	2.06	0.56
1:F:25:VAL:HG13	1:F:62:ALA:CB	2.35	0.56
1:D:83:HIS:CD2	1:D:83:HIS:H	2.24	0.56
1:F:168:PRO:HB2	1:F:169:GLU:OE1	2.05	0.55
1:E:122:GLN:N	1:E:122:GLN:HE21	1.97	0.55
1:A:250:ALA:O	1:A:254:ARG:HG3	2.06	0.55

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:28:LEU:HD23	1:A:116:LEU:HD21	1.88	0.55
1:F:28:LEU:HD23	1:F:116:LEU:HD11	1.89	0.55
1:C:176:ILE:N	1:C:176:ILE:HD13	2.22	0.55
1:F:28:LEU:HD23	1:F:116:LEU:CD1	2.37	0.54
1:D:151:LEU:O	1:D:151:LEU:HD12	2.07	0.54
1:B:28:LEU:HD23	1:B:116:LEU:HD21	1.88	0.53
1:E:235:ARG:O	1:E:239:GLU:HG3	2.09	0.53
1:A:235:ARG:O	1:A:239:GLU:HG3	2.09	0.53
1:C:78:SER:OG	1:C:122:GLN:NE2	2.42	0.53
1:D:172:LEU:HB2	1:D:176:ILE:HD11	1.90	0.53
1:B:39:HIS:CD2	1:B:40:MET:HG3	2.43	0.53
1:D:151:LEU:HD22	1:D:156:PHE:CE1	2.44	0.52
1:D:116:LEU:HD12	1:D:131:ILE:HD11	1.92	0.52
1:D:159:ASP:OD2	1:D:161:ALA:HB3	2.10	0.52
1:B:171:PHE:C	1:B:172:LEU:HD12	2.30	0.52
1:D:26:LEU:HD23	1:D:63:LEU:HB2	1.91	0.52
1:A:176:ILE:HG12	1:A:176:ILE:O	2.09	0.52
1:E:176:ILE:CD1	1:E:205:PRO:HB3	2.40	0.51
1:F:217:ASP:HB3	1:F:220:ARG:HG3	1.91	0.51
1:D:26:LEU:CD1	1:D:106:LEU:HD22	2.40	0.51
1:B:216:VAL:HG11	4:B:1359:HOH:O	2.10	0.51
1:F:56:VAL:O	1:F:60:ILE:HG12	2.10	0.51
1:E:151:LEU:HD12	1:E:151:LEU:C	2.31	0.51
1:F:214:PHE:HA	1:F:215:PRO:C	2.32	0.51
1:B:56:VAL:O	1:B:60:ILE:HG12	2.11	0.51
1:F:39:HIS:CD2	1:F:40:MET:HG3	2.46	0.50
1:C:131:ILE:HD11	1:C:150:VAL:HG21	1.92	0.50
1:A:213:VAL:HG21	1:A:220:ARG:NH2	2.27	0.50
1:D:235:ARG:HD2	4:D:1491:HOH:O	2.11	0.50
1:A:39:HIS:CD2	1:A:40:MET:HG3	2.47	0.49
1:B:28:LEU:HD23	1:B:116:LEU:CD2	2.43	0.49
1:A:165:GLN:NE2	1:A:240:LEU:HD22	2.27	0.49
1:B:172:LEU:N	1:B:172:LEU:HD12	2.28	0.49
1:C:121:TYR:HB3	1:C:122:GLN:NE2	2.28	0.49
1:F:76:GLN:HG3	1:F:221:THR:OG1	2.13	0.49
1:D:121:TYR:HB3	1:D:122:GLN:OE1	2.13	0.48
1:F:151:LEU:HD12	1:F:151:LEU:C	2.34	0.48
1:E:26:LEU:HD13	1:E:106:LEU:HD22	1.95	0.48
1:A:151:LEU:C	1:A:151:LEU:HD12	2.33	0.48
1:B:77:LYS:HE2	1:B:212:ASP:OD1	2.12	0.48
1:F:159:ASP:OD2	1:F:161:ALA:HB3	2.13	0.48
1:D:159:ASP:O	1:D:163:ILE:HD13	2.13	0.48

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:157:VAL:HG21	1:F:241:ILE:HG23	1.95	0.48
1:E:168:PRO:C	1:E:170:GLY:H	2.16	0.48
1:A:117:MET:HG3	1:A:151:LEU:HD11	1.96	0.48
1:A:150:VAL:CG1	1:B:212:ASP:HB2	2.43	0.48
1:C:55:ARG:HG3	1:C:55:ARG:HH11	1.79	0.47
1:C:242:LEU:O	1:C:246:VAL:HG23	2.15	0.47
1:B:12:TRP:CH2	1:B:13:LYS:HE3	2.49	0.47
1:A:150:VAL:HG13	1:B:212:ASP:HB2	1.96	0.47
1:B:3:LYS:NZ	4:B:1309:HOH:O	2.47	0.47
1:C:214:PHE:HA	1:C:215:PRO:C	2.33	0.47
1:D:214:PHE:HA	1:D:215:PRO:C	2.34	0.47
1:C:28:LEU:HD23	1:C:116:LEU:HD21	1.97	0.47
1:B:209:PRO:HB2	1:B:211:TYR:CE2	2.50	0.47
1:E:159:ASP:OD2	1:E:162:VAL:HG23	2.15	0.47
1:D:12:TRP:CD1	1:E:35:GLN:HB2	2.49	0.47
1:B:140:TYR:OH	1:F:137:GLU:HG2	2.15	0.47
1:D:116:LEU:HD13	1:D:127:ILE:HG23	1.96	0.46
1:F:25:VAL:HG13	1:F:62:ALA:HB2	1.96	0.46
1:D:54:LYS:O	1:D:58:GLU:HG3	2.16	0.46
1:F:120:HIS:CE1	1:F:122:GLN:HG2	2.51	0.46
1:C:173:GLY:C	1:C:175:ASP:H	2.19	0.46
1:A:175:ASP:OD1	1:A:176:ILE:N	2.49	0.46
1:D:55:ARG:HG3	1:D:55:ARG:HH11	1.80	0.45
1:E:159:ASP:O	1:E:163:ILE:HG12	2.16	0.45
1:F:176:ILE:C	1:F:176:ILE:HD12	2.36	0.45
1:A:209:PRO:HB2	1:A:211:TYR:CE2	2.51	0.45
1:F:26:LEU:HD13	1:F:106:LEU:HD22	1.99	0.45
1:C:131:ILE:HD11	1:C:150:VAL:CG2	2.46	0.45
1:E:214:PHE:HA	1:E:215:PRO:C	2.36	0.45
1:A:221:THR:HG23	4:A:1156:HOH:O	2.16	0.45
1:D:60:ILE:HD12	1:D:253:ILE:HG21	1.99	0.45
1:A:214:PHE:HA	1:A:215:PRO:C	2.37	0.45
1:E:175:ASP:OD2	1:E:176:ILE:N	2.50	0.45
1:E:122:GLN:N	1:E:122:GLN:NE2	2.61	0.44
1:D:116:LEU:HD13	1:D:127:ILE:CG2	2.48	0.44
1:B:73:LYS:O	1:B:122:GLN:HG3	2.17	0.44
4:C:1111:HOH:O	1:D:125:MET:HG3	2.16	0.44
1:E:26:LEU:HD23	1:E:63:LEU:HB2	1.99	0.44
1:E:171:PHE:CG	1:E:172:LEU:N	2.86	0.44
1:F:169:GLU:N	1:F:169:GLU:OE1	2.51	0.44
1:A:72:TYR:CG	1:A:73:LYS:N	2.85	0.43
1:B:17:ALA:O	1:B:20:ALA:HB3	2.18	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:18:ARG:NH1	4:D:1082:HOH:O	2.50	0.43
1:E:176:ILE:O	1:E:176:ILE:HG12	2.18	0.43
1:D:167:TYR:HB3	1:D:170:GLY:O	2.18	0.43
1:C:131:ILE:CD1	1:C:150:VAL:CG2	2.97	0.43
1:D:223:ALA:N	1:D:224:PRO:HD2	2.32	0.43
1:B:69:GLN:NE2	4:B:1418:HOH:O	2.52	0.43
1:C:209:PRO:HB2	1:C:211:TYR:CE2	2.54	0.43
1:E:197:LEU:O	1:E:200:VAL:HG23	2.18	0.43
1:B:159:ASP:OD1	1:B:161:ALA:N	2.52	0.43
1:E:153:TYR:HE1	1:E:182:PHE:HE1	1.67	0.43
1:E:72:TYR:CG	1:E:73:LYS:N	2.87	0.42
1:A:12:TRP:CE3	1:F:85:PRO:HD2	2.55	0.42
1:E:216:VAL:O	1:E:218:PRO:HD3	2.19	0.42
1:B:176:ILE:HG12	1:B:176:ILE:O	2.19	0.42
1:A:120:HIS:CE1	1:A:122:GLN:NE2	2.87	0.42
1:C:75:GLN:CG	1:C:122:GLN:HG2	2.33	0.42
1:E:242:LEU:O	1:E:246:VAL:HG23	2.20	0.42
1:D:35:GLN:HB2	1:E:12:TRP:CD1	2.55	0.42
1:A:45:ASP:O	1:A:49:PRO:HG3	2.20	0.42
1:B:72:TYR:CG	1:B:73:LYS:N	2.88	0.42
1:F:216:VAL:O	1:F:218:PRO:HD3	2.20	0.42
1:E:177:GLU:HG3	1:E:177:GLU:O	2.20	0.42
1:B:258:PRO:HA	1:B:259:PRO:HD3	1.91	0.42
1:D:151:LEU:HD12	1:D:151:LEU:C	2.40	0.42
1:A:176:ILE:HD11	1:A:205:PRO:HB3	2.01	0.42
1:A:26:LEU:CD1	1:A:106:LEU:HD22	2.49	0.42
1:F:169:GLU:N	1:F:169:GLU:CD	2.73	0.42
1:C:48:LEU:O	1:C:52:VAL:HG23	2.20	0.42
1:B:214:PHE:HA	1:B:215:PRO:C	2.39	0.42
1:A:79:GLY:HA3	4:A:1157:HOH:O	2.19	0.42
1:E:196:ASP:OD2	1:E:199:ARG:NH2	2.53	0.42
1:A:28:LEU:HD23	1:A:116:LEU:CD2	2.50	0.41
1:D:100:GLN:NE2	1:D:137:GLU:OE2	2.53	0.41
1:B:30:VAL:HG13	1:B:68:LEU:HD12	2.02	0.41
1:E:77:LYS:HD3	1:E:212:ASP:OD1	2.20	0.41
1:A:54:LYS:O	1:A:58:GLU:HG3	2.19	0.41
1:E:249:ILE:O	1:E:253:ILE:HG13	2.20	0.41
1:E:209:PRO:HB2	1:E:211:TYR:CE2	2.55	0.41
1:E:151:LEU:HD12	1:E:151:LEU:O	2.20	0.41
1:F:72:TYR:CG	1:F:73:LYS:N	2.88	0.41
1:D:85:PRO:HD2	1:E:12:TRP:CE3	2.55	0.41
1:E:129:GLU:O	1:E:133:LEU:HG	2.21	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:176:ILE:HG21	1:C:206:ALA:O	2.21	0.41
1:E:162:VAL:O	1:E:166:LEU:HG	2.20	0.41
1:F:151:LEU:O	1:F:151:LEU:HD12	2.21	0.41
1:E:56:VAL:O	1:E:60:ILE:HG12	2.20	0.41
1:A:117:MET:HG3	1:A:151:LEU:CD1	2.51	0.41
1:E:258:PRO:HA	1:E:259:PRO:HD3	1.83	0.41
1:B:18:ARG:HA	1:B:18:ARG:HD2	1.90	0.41
1:A:258:PRO:HA	1:A:259:PRO:HD3	1.89	0.41
1:A:65:MET:CE	1:A:106:LEU:HD21	2.51	0.41
1:F:234:SER:OG	1:F:237:LYS:HG3	2.20	0.41
1:E:117:MET:HG3	1:E:151:LEU:HD11	2.03	0.40
1:A:242:LEU:O	1:A:246:VAL:HG23	2.22	0.40
1:A:56:VAL:O	1:A:60:ILE:HG12	2.21	0.40
1:D:258:PRO:HA	1:D:259:PRO:HD3	1.85	0.40
1:F:129:GLU:HA	1:F:129:GLU:OE1	2.20	0.40
1:C:223:ALA:N	1:C:224:PRO:CD	2.85	0.40
1:F:139:ARG:HG3	1:F:144:GLN:HE21	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%)	243 (95%)	12 (5%)	0	100	100
1	B	255/260 (98%)	239 (94%)	16 (6%)	0	100	100
1	C	256/260 (98%)	242 (94%)	14 (6%)	0	100	100
1	D	254/260 (98%)	240 (94%)	13 (5%)	1 (0%)	43	36
1	E	254/260 (98%)	237 (93%)	16 (6%)	1 (0%)	43	36
1	F	255/260 (98%)	244 (96%)	11 (4%)	0	100	100
All	All	1529/1560 (98%)	1445 (94%)	82 (5%)	2 (0%)	59	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	169	GLU
1	D	173	GLY

### 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	208/215 (97%)	206 (99%)	2 (1%)	85	88
1	B	212/215 (99%)	209 (99%)	3 (1%)	78	81
1	C	212/215 (99%)	209 (99%)	3 (1%)	78	81
1	D	211/215 (98%)	211 (100%)	0	100	100
1	E	210/215 (98%)	207 (99%)	3 (1%)	78	81
1	F	212/215 (99%)	209 (99%)	3 (1%)	78	81
All	All	1265/1290 (98%)	1251 (99%)	14 (1%)	84	86

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

Mol	Chain	Res	Type
1	A	122	GLN
1	A	212	ASP
1	B	122	GLN
1	B	158	LYS
1	B	193	ASP
1	C	176	ILE
1	C	202	ASP
1	C	231	LYS
1	E	116	LEU
1	E	122	GLN
1	E	212	ASP
1	F	122	GLN
1	F	169	GLU
1	F	175	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	122	GLN
1	A	165	GLN
1	B	69	GLN
1	B	122	GLN
1	B	144	GLN
1	C	122	GLN
1	D	83	HIS
1	D	165	GLN
1	E	76	GLN
1	E	82	ASN
1	E	122	GLN
1	F	76	GLN
1	F	122	GLN
1	F	144	GLN
1	F	165	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 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.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	257/260 (98%)	-0.12	6 (2%) 57 57	17, 26, 47, 61	0
1	B	257/260 (98%)	-0.15	7 (2%) 52 52	16, 26, 49, 64	0
1	C	258/260 (99%)	-0.14	9 (3%) 42 41	16, 28, 51, 63	0
1	D	256/260 (98%)	-0.10	11 (4%) 34 32	15, 28, 50, 62	0
1	E	256/260 (98%)	-0.17	8 (3%) 47 46	15, 28, 49, 58	0
1	F	257/260 (98%)	-0.23	4 (1%) 68 69	15, 28, 49, 62	0
All	All	1541/1560 (98%)	-0.15	45 (2%) 49 49	15, 28, 50, 64	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	170	GLY	5.0
1	D	172	LEU	4.4
1	B	171	PHE	4.2
1	D	171	PHE	4.2
1	E	172	LEU	4.1
1	D	170	GLY	3.9
1	C	175	ASP	3.9
1	E	22	GLY	3.7
1	A	175	ASP	3.6
1	B	169	GLU	3.5
1	C	170	GLY	3.5
1	D	160	PRO	3.3
1	F	169	GLU	3.2
1	E	259	PRO	3.1
1	D	169	GLU	3.1
1	C	172	LEU	3.0
1	C	202	ASP	3.0
1	C	171	PHE	3.0
1	B	175	ASP	2.9

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	172	LEU	2.9
1	C	2	SER	2.9
1	E	169	GLU	2.9
1	A	22	GLY	2.8
1	A	171	PHE	2.8
1	E	170	GLY	2.7
1	F	164	GLN	2.7
1	A	168	PRO	2.7
1	C	160	PRO	2.6
1	C	164	GLN	2.6
1	D	20	ALA	2.6
1	D	175	ASP	2.6
1	F	172	LEU	2.5
1	B	168	PRO	2.5
1	A	206	ALA	2.5
1	D	22	GLY	2.4
1	D	176	ILE	2.4
1	E	164	GLN	2.4
1	C	165	GLN	2.2
1	D	259	PRO	2.2
1	D	174	TRP	2.2
1	E	174	TRP	2.2
1	E	176	ILE	2.1
1	F	207	THR	2.1
1	B	198	ASP	2.1
1	A	121	TYR	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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	CL	C	302	1/1	0.28	6.62	39,39,39,39	0
3	CL	F	302	1/1	0.25	5.89	34,34,34,34	0
3	CL	B	302	1/1	0.26	5.03	34,34,34,34	0
3	CL	D	302	1/1	0.27	4.65	34,34,34,34	0
3	CL	E	302	1/1	0.26	4.43	33,33,33,33	0
3	CL	A	302	1/1	0.28	3.80	36,36,36,36	0
2	ZN	D	301	1/1	0.13	1.11	25,25,25,25	0
2	ZN	E	301	1/1	0.10	0.10	25,25,25,25	0
2	ZN	B	301	1/1	0.10	0.06	23,23,23,23	0
2	ZN	F	301	1/1	0.10	-0.04	24,24,24,24	0
2	ZN	A	301	1/1	0.09	-0.61	24,24,24,24	0
2	ZN	E	300	1/1	0.09	-0.67	26,26,26,26	0
2	ZN	C	301	1/1	0.09	-0.75	26,26,26,26	0
2	ZN	C	300	1/1	0.09	-0.80	29,29,29,29	0
2	ZN	B	300	1/1	0.08	-1.07	25,25,25,25	0
2	ZN	F	300	1/1	0.07	-1.23	27,27,27,27	0
2	ZN	A	300	1/1	0.07	-1.37	26,26,26,26	0
2	ZN	D	300	1/1	0.06	-2.08	25,25,25,25	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.