Full wwPDB X-ray Structure Validation Report (i)

Dec 19, 2023 – 04:14 PM EST

PDB ID	:	1J2T
Title	:	Creatininase Mn
Authors	:	Yoshimoto, T.; Tanaka, N.; Kanada, N.; Inoue, T.; Nakajima, Y.; Haratake, M.; Nakamura, K.T.; Xu, Y.; Ito, K.
Deposited on	:	2003-01-11
Resolution	:	1.80 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

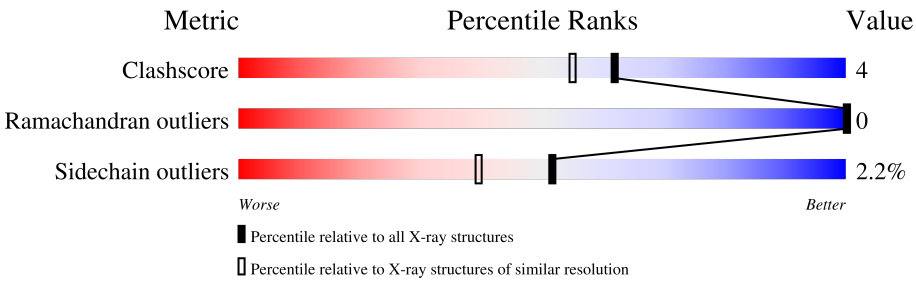
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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))
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	260	90% 8% ••
1	B	260	85% 11% ••
1	C	260	88% 10% ••
1	D	260	89% 9% ••
1	E	260	87% 11% ••
1	F	260	90% 8% ••

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 13386 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
			Total	C	N	O	S			
1	A	257	1988	1273	337	367	11	0	0	0
1	B	257	1988	1273	337	367	11	0	0	0
1	C	257	1988	1273	337	367	11	0	0	0
1	D	257	1988	1273	337	367	11	0	0	0
1	E	257	1988	1273	337	367	11	0	0	0
1	F	257	1988	1273	337	367	11	0	0	0

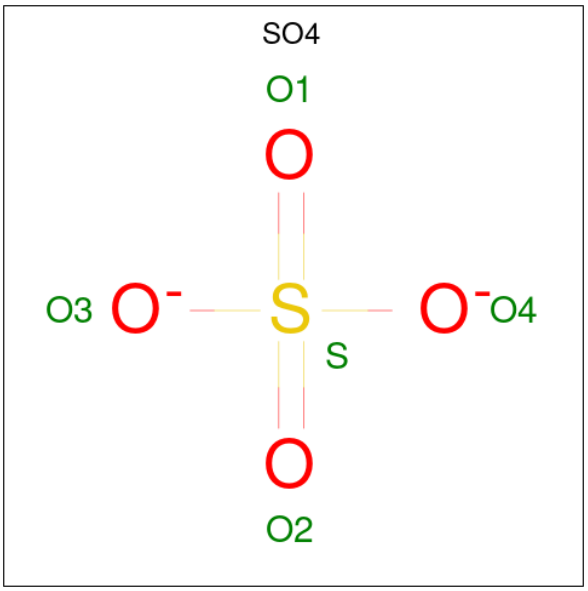
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
4	D	1	5	4	1	0	0
			Total	O	S		
4	E	1	5	4	1	0	0
			Total	O	S		
4	F	1	5	4	1	0	0
			Total	O	S		
4	F	1	5	4	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	267	267	267	0	0
			Total	O		
5	B	235	235	235	0	0
			Total	O		
5	C	214	214	214	0	0
			Total	O		
5	D	229	229	229	0	0
			Total	O		
5	E	204	204	204	0	0
			Total	O		
5	F	247	247	247	0	0

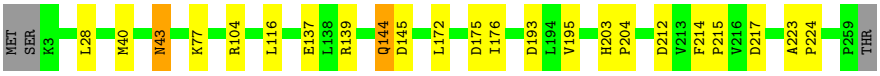
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

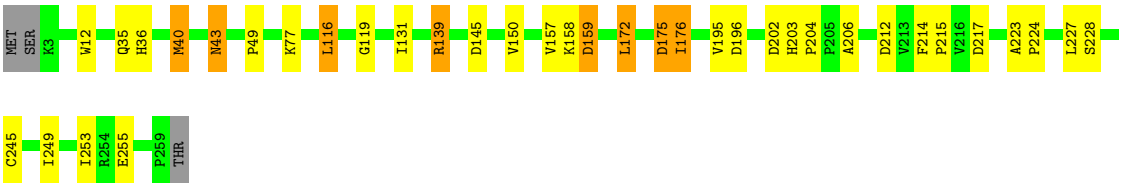
• Molecule 1: creatinine amidohydrolase

Chain A:



• Molecule 1: creatinine amidohydrolase

Chain B:



• Molecule 1: creatinine amidohydrolase

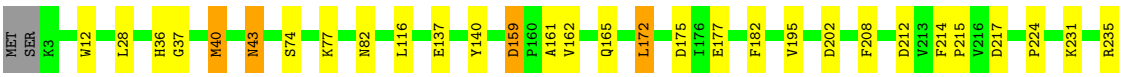
Chain C:

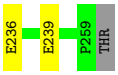
• Molecule 1: creatinine amidohydrolase

Chain D:

• Molecule 1: creatinine amidohydrolase

Chain E:





• Molecule 1: creatinine amidohydrolase

Chain F:

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	101.99Å 150.74Å 167.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.80	Depositor
% Data completeness (in resolution range)	99.9 (40.00-1.80)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.175 , 0.193	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	13386	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

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: MN, ZN, SO4

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	0.44	0/2037	0.75	4/2772 (0.1%)
1	B	0.42	0/2037	0.72	6/2772 (0.2%)
1	C	0.44	0/2037	0.85	9/2772 (0.3%)
1	D	0.43	0/2037	0.73	4/2772 (0.1%)
1	E	0.43	0/2037	0.73	3/2772 (0.1%)
1	F	0.43	0/2037	0.81	11/2772 (0.4%)
All	All	0.43	0/12222	0.76	37/16632 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	139	ARG	NE-CZ-NH1	-15.36	112.62	120.30
1	F	139	ARG	NE-CZ-NH1	-9.24	115.68	120.30
1	C	139	ARG	NE-CZ-NH2	8.29	124.45	120.30
1	C	139	ARG	CG-CD-NE	8.18	128.97	111.80
1	C	139	ARG	CD-NE-CZ	7.57	134.20	123.60
1	A	175	ASP	CB-CG-OD2	7.40	124.96	118.30
1	F	175	ASP	CB-CG-OD2	7.01	124.61	118.30
1	C	23	ASP	CB-CG-OD2	6.90	124.51	118.30
1	E	175	ASP	CB-CG-OD2	6.82	124.44	118.30
1	F	139	ARG	CG-CD-NE	6.76	126.00	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	F	139	ARG	NE-CZ-NH2	6.75	123.68	120.30
1	F	235	ARG	NE-CZ-NH1	-6.63	116.98	120.30
1	F	235	ARG	CG-CD-NE	6.52	125.49	111.80
1	F	235	ARG	NE-CZ-NH2	6.46	123.53	120.30
1	A	145	ASP	CB-CG-OD2	6.31	123.98	118.30
1	B	175	ASP	CB-CG-OD2	6.05	123.75	118.30
1	F	145	ASP	CB-CG-OD2	5.95	123.66	118.30
1	E	217	ASP	CB-CG-OD2	5.89	123.61	118.30
1	C	193	ASP	CB-CG-OD2	5.81	123.53	118.30
1	A	193	ASP	CB-CG-OD2	5.73	123.45	118.30
1	B	145	ASP	CB-CG-OD2	5.57	123.31	118.30
1	F	193	ASP	CB-CG-OD2	5.56	123.30	118.30
1	D	91	ASP	CB-CG-OD2	5.54	123.28	118.30
1	F	91	ASP	CB-CG-OD2	5.50	123.25	118.30
1	C	145	ASP	CB-CG-OD2	5.49	123.24	118.30
1	D	145	ASP	CB-CG-OD2	5.46	123.22	118.30
1	B	159	ASP	CB-CG-OD2	5.42	123.18	118.30
1	C	155	ASP	CB-CG-OD2	5.42	123.17	118.30
1	A	217	ASP	CB-CG-OD2	5.41	123.17	118.30
1	D	175	ASP	CB-CG-OD2	5.39	123.15	118.30
1	D	198	ASP	CB-CG-OD2	5.36	123.13	118.30
1	B	217	ASP	CB-CG-OD2	5.29	123.06	118.30
1	F	159	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	196	ASP	CB-CG-OD2	5.20	122.98	118.30
1	C	217	ASP	CB-CG-OD2	5.18	122.97	118.30
1	B	202	ASP	CB-CG-OD2	5.16	122.94	118.30
1	E	159	ASP	CB-CG-OD2	5.13	122.92	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	139	ARG	Sidechain

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	1988	0	1969	17	0
1	B	1988	0	1969	23	0
1	C	1988	0	1969	18	0
1	D	1988	0	1969	15	0
1	E	1988	0	1969	26	0
1	F	1988	0	1969	12	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	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	10	0	0	0	0
4	B	10	0	0	0	0
4	C	10	0	0	0	0
4	D	5	0	0	0	0
4	E	5	0	0	0	0
4	F	10	0	0	0	0
5	A	267	0	0	2	1
5	B	235	0	0	5	1
5	C	214	0	0	3	0
5	D	229	0	0	1	0
5	E	204	0	0	2	0
5	F	247	0	0	2	0
All	All	13386	0	11814	104	1

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 4.

All (104) 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:C:40:MET:CE	1:C:40:MET:SD	2.04	1.43
1:D:235:ARG:CZ	1:D:236:GLU:OE1	2.08	1.00
1:E:235:ARG:CZ	1:E:236:GLU:OE1	2.10	0.99
1:E:172:LEU:HD22	1:E:172:LEU:H	1.33	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:235:ARG:NH1	1:E:236:GLU:OE1	2.02	0.91
1:D:235:ARG:NH1	1:D:236:GLU:OE1	2.05	0.89
1:C:216:VAL:HG12	5:C:7128:HOH:O	1.76	0.85
1:F:13:LYS:NZ	5:F:7201:HOH:O	2.09	0.79
1:D:77:LYS:HG2	5:D:7059:HOH:O	1.84	0.77
1:B:175:ASP:HB3	5:B:7235:HOH:O	1.91	0.70
1:E:236:GLU:OE2	5:E:7142:HOH:O	2.09	0.69
1:C:175:ASP:HB3	5:C:7216:HOH:O	1.92	0.69
1:E:172:LEU:HD22	1:E:172:LEU:N	2.09	0.67
1:E:177:GLU:CD	1:E:182:PHE:HB3	2.19	0.62
1:B:157:VAL:CG2	1:B:245:CYS:SG	2.88	0.62
1:C:140:TYR:OH	1:E:137:GLU:OE1	2.18	0.62
1:B:176:ILE:HD11	1:B:206:ALA:O	2.01	0.61
1:B:139:ARG:NE	5:B:7098:HOH:O	1.96	0.59
1:E:161:ALA:O	1:E:165:GLN:HG3	2.03	0.59
1:A:104:ARG:HD2	5:A:7205:HOH:O	2.03	0.58
1:B:40:MET:HG3	1:B:195:VAL:HG22	1.84	0.58
1:F:40:MET:HG3	1:F:195:VAL:HG22	1.85	0.58
1:E:36:HIS:HB3	1:E:40:MET:HE3	1.86	0.58
1:C:214:PHE:HA	1:C:215:PRO:C	2.25	0.57
1:D:159:ASP:OD2	1:D:161:ALA:HB3	2.07	0.54
1:B:131:ILE:CD1	1:B:150:VAL:HG21	2.38	0.53
1:B:172:LEU:HB2	5:B:7159:HOH:O	2.08	0.53
1:E:235:ARG:NH2	1:E:236:GLU:OE1	2.41	0.53
1:D:40:MET:HG3	1:D:195:VAL:HG22	1.92	0.52
1:B:77:LYS:HD2	1:B:212:ASP:OD2	2.10	0.52
1:F:214:PHE:HA	1:F:215:PRO:C	2.29	0.52
1:B:43:ASN:C	1:B:43:ASN:HD22	2.13	0.51
1:D:214:PHE:HA	1:D:215:PRO:C	2.30	0.51
1:B:214:PHE:HA	1:B:215:PRO:C	2.31	0.50
1:A:28:LEU:HD23	1:A:116:LEU:HD21	1.93	0.50
1:A:214:PHE:HA	1:A:215:PRO:C	2.31	0.50
1:B:176:ILE:HG12	1:B:227:LEU:O	2.12	0.49
1:C:175:ASP:CB	5:C:7216:HOH:O	2.58	0.49
1:E:214:PHE:HA	1:E:215:PRO:C	2.31	0.49
1:A:77:LYS:CD	1:A:212:ASP:OD2	2.61	0.49
1:F:74:SER:H	1:F:82:ASN:HD22	1.60	0.49
1:C:40:MET:HG3	1:C:195:VAL:HG22	1.95	0.48
1:B:36:HIS:HB3	1:B:40:MET:HE3	1.96	0.48
1:D:187:MET:HB3	1:D:195:VAL:HG21	1.96	0.48
1:A:77:LYS:NZ	5:A:7272:HOH:O	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ARG:NH2	1:A:144:GLN:OE1	2.48	0.47
1:C:131:ILE:CD1	1:C:150:VAL:HG21	2.44	0.47
1:E:28:LEU:HD23	1:E:116:LEU:HD21	1.97	0.47
1:A:43:ASN:HD22	1:A:43:ASN:C	2.18	0.47
1:C:43:ASN:C	1:C:43:ASN:HD22	2.18	0.47
1:A:77:LYS:HD2	1:A:212:ASP:OD2	2.15	0.47
1:D:43:ASN:HD22	1:D:43:ASN:C	2.19	0.46
1:F:158:LYS:NZ	5:F:7242:HOH:O	2.30	0.46
1:B:49:PRO:HG3	1:B:119:GLY:HA3	1.99	0.45
1:F:37:GLY:HA2	1:F:224:PRO:O	2.16	0.45
1:F:43:ASN:C	1:F:43:ASN:HD22	2.19	0.45
1:A:40:MET:HG3	1:A:195:VAL:HG22	1.99	0.45
1:E:212:ASP:HB2	1:F:150:VAL:CG1	2.47	0.45
1:C:40:MET:HE2	1:C:40:MET:CB	2.47	0.45
1:D:74:SER:H	1:D:82:ASN:HD22	1.63	0.45
1:A:223:ALA:HB3	1:A:224:PRO:HD3	1.98	0.45
1:B:158:LYS:HG3	5:B:7173:HOH:O	2.18	0.44
1:E:43:ASN:HD22	1:E:43:ASN:C	2.20	0.44
1:B:159:ASP:C	1:B:159:ASP:OD1	2.56	0.44
1:B:176:ILE:HG23	1:B:228:SER:HB3	1.99	0.44
1:A:176:ILE:CG2	1:A:176:ILE:O	2.65	0.44
1:B:249:ILE:O	1:B:253:ILE:HG12	2.18	0.44
1:E:77:LYS:HG2	1:E:208:PHE:CD1	2.53	0.44
1:D:187:MET:HB3	1:D:195:VAL:CG2	2.48	0.43
1:C:159:ASP:HA	1:C:160:PRO:HD2	1.92	0.43
1:F:36:HIS:HB3	1:F:40:MET:HE3	2.00	0.43
1:F:223:ALA:HB3	1:F:224:PRO:HD3	2.01	0.43
1:B:12:TRP:CD1	1:C:35:GLN:HB2	2.53	0.43
1:C:40:MET:CE	1:C:40:MET:CG	2.94	0.43
1:E:37:GLY:HA2	1:E:224:PRO:O	2.19	0.43
1:E:159:ASP:OD2	1:E:161:ALA:HB3	2.19	0.43
1:D:26:LEU:CD1	1:D:106:LEU:HD22	2.49	0.42
1:B:116:LEU:HD13	1:B:131:ILE:HD11	2.01	0.42
1:D:223:ALA:HB3	1:D:224:PRO:HD3	2.01	0.42
1:A:139:ARG:CZ	1:A:144:GLN:OE1	2.67	0.42
1:E:40:MET:HG3	1:E:195:VAL:HG22	1.99	0.42
1:A:144:GLN:HE21	1:A:144:GLN:HB2	1.57	0.42
1:A:203:HIS:HB2	1:A:204:PRO:HD2	2.01	0.42
1:E:159:ASP:OD2	1:E:162:VAL:HG23	2.20	0.42
1:E:202:ASP:CG	1:E:231:LYS:HD2	2.39	0.42
1:C:223:ALA:N	1:C:224:PRO:CD	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:ALA:N	1:B:224:PRO:CD	2.83	0.42
1:C:37:GLY:HA2	1:C:224:PRO:O	2.19	0.41
1:D:236:GLU:H	1:D:236:GLU:CD	2.24	0.41
1:A:137:GLU:OE1	1:E:140:TYR:OH	2.38	0.41
1:A:176:ILE:O	1:A:176:ILE:HG22	2.21	0.41
1:F:116:LEU:HD22	1:F:127:ILE:HG23	2.03	0.41
1:C:74:SER:H	1:C:82:ASN:HD22	1.68	0.41
1:D:35:GLN:HB2	1:E:12:TRP:CD1	2.56	0.41
1:A:77:LYS:HD3	1:A:212:ASP:OD2	2.21	0.41
1:B:139:ARG:HG3	5:B:7098:HOH:O	2.21	0.41
1:B:203:HIS:HB2	1:B:204:PRO:HD2	2.03	0.40
1:D:139:ARG:NH2	1:D:144:GLN:OE1	2.55	0.40
1:E:74:SER:H	1:E:82:ASN:HD22	1.69	0.40
1:E:172:LEU:N	1:E:172:LEU:CD2	2.78	0.40
1:B:35:GLN:HB2	1:C:12:TRP:CD1	2.57	0.40
1:C:40:MET:CE	1:C:40:MET:CB	3.00	0.40
1:E:212:ASP:HB2	1:F:150:VAL:HG13	2.03	0.40
1:E:239:GLU:HG2	5:E:7185:HOH:O	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:7197:HOH:O	5:B:7177:HOH:O[4_555]	2.18	0.02

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%)	251 (98%)	4 (2%)	0	100	100
1	B	255/260 (98%)	250 (98%)	5 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	255/260 (98%)	250 (98%)	5 (2%)	0	100	100
1	D	255/260 (98%)	251 (98%)	4 (2%)	0	100	100
1	E	255/260 (98%)	251 (98%)	4 (2%)	0	100	100
1	F	255/260 (98%)	252 (99%)	3 (1%)	0	100	100
All	All	1530/1560 (98%)	1505 (98%)	25 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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	212/215 (99%)	209 (99%)	3 (1%)	67	59
1	B	212/215 (99%)	205 (97%)	7 (3%)	38	23
1	C	212/215 (99%)	206 (97%)	6 (3%)	43	30
1	D	212/215 (99%)	208 (98%)	4 (2%)	57	46
1	E	212/215 (99%)	209 (99%)	3 (1%)	67	59
1	F	212/215 (99%)	207 (98%)	5 (2%)	49	36
All	All	1272/1290 (99%)	1244 (98%)	28 (2%)	52	39

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	144	GLN
1	A	172	LEU
1	B	40	MET
1	B	43	ASN
1	B	116	LEU
1	B	139	ARG
1	B	172	LEU
1	B	176	ILE

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Mol	Chain	Res	Type
1	B	255	GLU
1	C	40	MET
1	C	43	ASN
1	C	116	LEU
1	C	139	ARG
1	C	172	LEU
1	C	255	GLU
1	D	40	MET
1	D	43	ASN
1	D	172	LEU
1	D	177	GLU
1	E	40	MET
1	E	43	ASN
1	E	172	LEU
1	F	3	LYS
1	F	40	MET
1	F	43	ASN
1	F	116	LEU
1	F	139	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	82	ASN
1	B	43	ASN
1	B	82	ASN
1	C	43	ASN
1	C	75	GLN
1	C	82	ASN
1	D	43	ASN
1	D	75	GLN
1	D	82	ASN
1	E	43	ASN
1	E	82	ASN
1	E	144	GLN
1	F	43	ASN
1	F	75	GLN
1	F	82	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 22 ligands modelled in this entry, 12 are monoatomic - leaving 10 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	SO4	E	7003	-	4,4,4	0.15	0	6,6,6	0.35	0
4	SO4	B	7002	-	4,4,4	0.20	0	6,6,6	0.17	0
4	SO4	C	7009	-	4,4,4	0.16	0	6,6,6	0.13	0
4	SO4	A	7007	-	4,4,4	0.18	0	6,6,6	0.15	0
4	SO4	F	7004	-	4,4,4	0.17	0	6,6,6	0.29	0
4	SO4	D	7001	-	4,4,4	0.13	0	6,6,6	0.23	0
4	SO4	A	7006	-	4,4,4	0.22	0	6,6,6	0.46	0
4	SO4	B	7008	-	4,4,4	0.14	0	6,6,6	0.06	0
4	SO4	C	7005	-	4,4,4	0.17	0	6,6,6	0.29	0
4	SO4	F	7010	-	4,4,4	0.13	0	6,6,6	0.16	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.

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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.