Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 08:07 pm BST

PDB ID	:	1O8U
Title	:	The 2 Angstrom Structure of 6-Oxo Camphor Hydrolase: New Structural Diversity in the Crotonase Superfamily
Authors	:	Grogan, G.; Whittingham, J.L.; Turkenburg, J.P.; Verma, C.S.; Walsh, M.A.
Deposited on	:	2002-12-04
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
https://www.wwpdb.org/validation/2017/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:

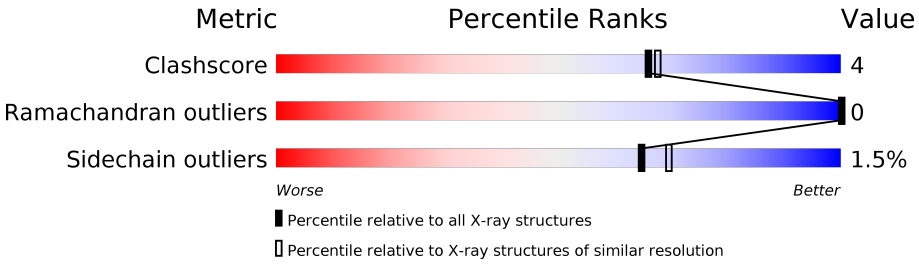
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 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))
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (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 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	257	89% 7% •
1	B	257	84% 13% •
1	C	257	86% 10% •
1	D	257	86% 9% • •
1	E	257	86% 9% • •
1	F	257	91% 6% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12473 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 6-OXO CAMPHOR HYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	0	0
			1910	1215	331	359	5			
1	B	249	Total	C	N	O	S	0	0	0
			1922	1222	332	363	5			
1	C	247	Total	C	N	O	S	0	1	0
			1913	1214	334	360	5			
1	D	247	Total	C	N	O	S	0	3	0
			1894	1208	323	358	5			
1	E	249	Total	C	N	O	S	0	1	0
			1926	1224	334	363	5			
1	F	248	Total	C	N	O	S	0	0	0
			1897	1208	323	361	5			

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		
2	D	1	Total	Na	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	179	Total	O	0	2
			179	179		
3	B	174	Total	O	0	0
			174	174		
3	C	170	Total	O	0	0
			170	170		
3	D	147	Total	O	0	0
			147	147		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	179	Total O 179 179	0	0
3	F	160	Total O 160 160	0	0

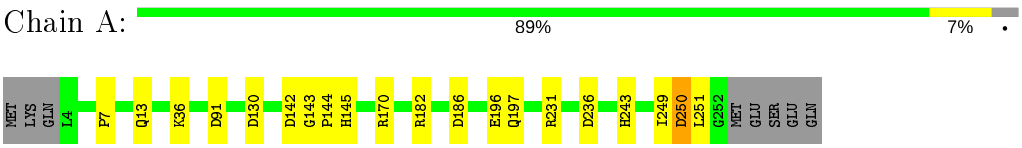
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

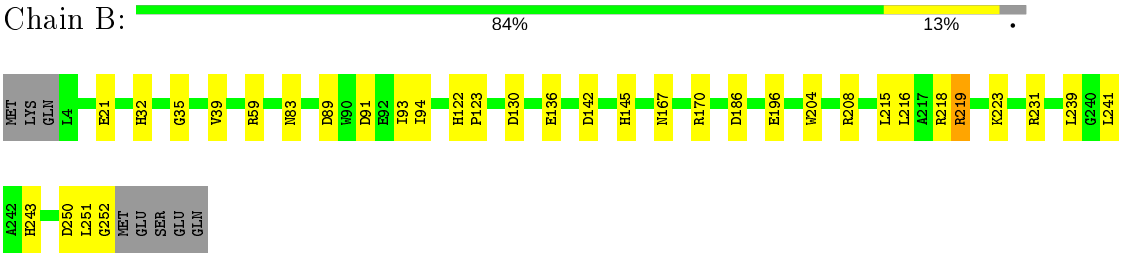
- Molecule 1: 6-OXO CAMPHOR HYDROLASE

Chain A:



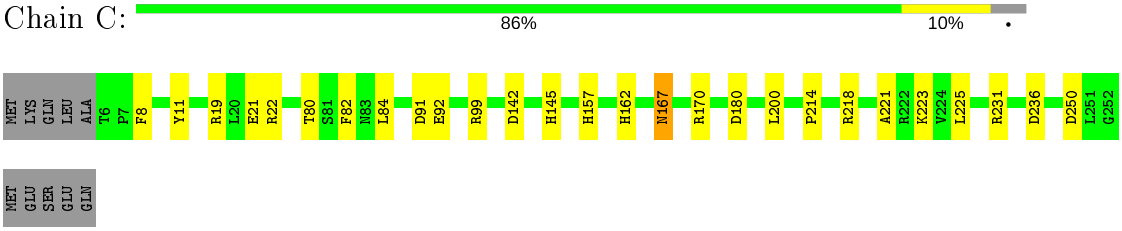
- Molecule 1: 6-OXO CAMPHOR HYDROLASE

Chain B:



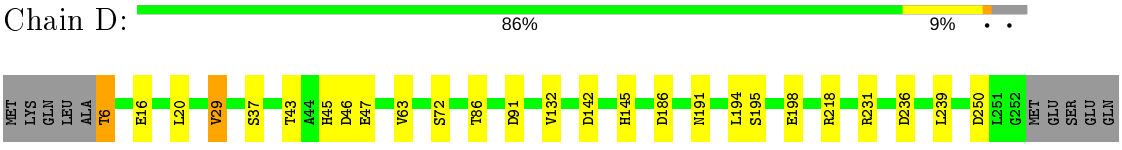
- Molecule 1: 6-OXO CAMPHOR HYDROLASE

Chain C:



- Molecule 1: 6-OXO CAMPHOR HYDROLASE

Chain D:

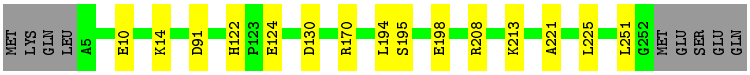


- Molecule 1: 6-OXO CAMPHOR HYDROLASE

Chain E:

- Molecule 1: 6-OXO CAMPHOR HYDROLASE

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 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	78.95Å 130.41Å 81.32Å 90.00° 114.16° 90.00°	Depositor
Resolution (Å)	25.00 – 2.00	Depositor
% Data completeness (in resolution range)	99.6 (25.00-2.00)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.147 , 0.190	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	12473	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA

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	0.86	4/1954 (0.2%)	1.12	12/2667 (0.4%)
1	B	0.85	0/1966	0.97	9/2684 (0.3%)
1	C	0.81	0/1964	1.01	9/2679 (0.3%)
1	D	0.79	0/1953	1.02	10/2668 (0.4%)
1	E	0.81	1/1975 (0.1%)	0.96	9/2695 (0.3%)
1	F	0.75	0/1940	0.89	5/2650 (0.2%)
All	All	0.81	5/11752 (0.0%)	1.00	54/16043 (0.3%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	249	ILE	C-O	9.70	1.41	1.23
1	A	251	LEU	CA-CB	6.84	1.69	1.53
1	A	249	ILE	C-N	-5.98	1.20	1.34
1	E	218	ARG	CB-CG	-5.18	1.38	1.52
1	A	196	GLU	CD-OE2	5.05	1.31	1.25

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	250	ASP	CB-CG-OD1	-19.25	100.97	118.30
1	A	231	ARG	NE-CZ-NH2	-18.62	110.99	120.30
1	A	250	ASP	CB-CG-OD2	18.60	135.04	118.30
1	D	231	ARG	NE-CZ-NH1	16.78	128.69	120.30
1	D	231	ARG	NE-CZ-NH2	-15.83	112.39	120.30
1	C	22	ARG	NE-CZ-NH2	-12.77	113.92	120.30
1	A	231	ARG	NE-CZ-NH1	12.32	126.46	120.30
1	C	22	ARG	NE-CZ-NH1	11.81	126.20	120.30
1	B	231	ARG	NE-CZ-NH1	11.60	126.10	120.30
1	B	231	ARG	NE-CZ-NH2	-11.41	114.59	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	231	ARG	NE-CZ-NH1	11.03	125.81	120.30
1	E	231	ARG	NE-CZ-NH1	9.45	125.03	120.30
1	C	231	ARG	NE-CZ-NH2	-9.44	115.58	120.30
1	D	186	ASP	CB-CG-OD2	8.56	126.00	118.30
1	C	91	ASP	CB-CG-OD2	8.20	125.68	118.30
1	E	250	ASP	CB-CG-OD2	8.00	125.50	118.30
1	E	208[A]	ARG	NE-CZ-NH1	7.53	124.07	120.30
1	E	208[B]	ARG	NE-CZ-NH1	7.53	124.07	120.30
1	A	91	ASP	CB-CG-OD2	7.48	125.03	118.30
1	F	130	ASP	CB-CG-OD2	7.45	125.01	118.30
1	E	231	ARG	NE-CZ-NH2	-6.81	116.89	120.30
1	E	218	ARG	NE-CZ-NH2	-6.65	116.97	120.30
1	E	99	ARG	NE-CZ-NH2	-6.45	117.08	120.30
1	B	130	ASP	CB-CG-OD2	6.27	123.94	118.30
1	A	231	ARG	CG-CD-NE	-6.21	98.77	111.80
1	E	130	ASP	CB-CG-OD2	5.87	123.58	118.30
1	F	170	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	B	91	ASP	CB-CG-OD1	5.83	123.55	118.30
1	A	231	ARG	CD-NE-CZ	5.82	131.74	123.60
1	A	186	ASP	CB-CG-OD2	5.82	123.53	118.30
1	F	208	ARG	NE-CZ-NH2	-5.73	117.43	120.30
1	C	218[A]	ARG	NE-CZ-NH1	5.69	123.15	120.30
1	C	218[B]	ARG	NE-CZ-NH1	5.69	123.15	120.30
1	B	186	ASP	CB-CG-OD2	5.57	123.31	118.30
1	A	170	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	D	231	ARG	CD-NE-CZ	5.55	131.37	123.60
1	A	251	LEU	CB-CA-C	-5.48	99.79	110.20
1	F	194	LEU	CA-CB-CG	5.47	127.88	115.30
1	B	219	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	F	91	ASP	CB-CG-OD2	5.37	123.13	118.30
1	D	46	ASP	CB-CG-OD2	5.34	123.10	118.30
1	E	99	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	A	236	ASP	CB-CG-OD2	5.30	123.07	118.30
1	C	180	ASP	CB-CG-OD1	5.25	123.02	118.30
1	B	142	ASP	CB-CG-OD2	5.24	123.02	118.30
1	C	250	ASP	CB-CG-OD2	5.24	123.01	118.30
1	B	231	ARG	CD-NE-CZ	5.14	130.80	123.60
1	D	91	ASP	CB-CG-OD2	5.10	122.89	118.30
1	B	231	ARG	CA-CB-CG	5.09	124.61	113.40
1	A	130	ASP	CB-CG-OD2	5.09	122.88	118.30
1	D	194	LEU	CA-CB-CG	5.07	126.96	115.30
1	D	250	ASP	CB-CG-OD2	5.04	122.83	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	236	ASP	CB-CG-OD2	5.01	122.81	118.30
1	D	231	ARG	CA-CB-CG	5.00	124.41	113.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	1910	0	1857	12	0
1	B	1922	0	1872	30	0
1	C	1913	0	1862	15	0
1	D	1894	0	1837	11	0
1	E	1926	0	1884	27	0
1	F	1897	0	1839	7	0
2	A	1	0	0	0	0
2	D	1	0	0	0	0
3	A	179	0	0	5	0
3	B	174	0	0	7	1
3	C	170	0	0	4	0
3	D	147	0	0	2	1
3	E	179	0	0	6	0
3	F	160	0	0	3	0
All	All	12473	0	11151	91	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 (91) 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:195:SER:OG	1:F:198:GLU:HG3	1.71	0.89
1:B:215:LEU:HD11	1:B:219:ARG:CZ	2.04	0.87
1:D:195:SER:OG	1:D:198:GLU:HG3	1.78	0.83
1:B:32:HIS:HD2	1:B:35:GLY:H	1.25	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:21:GLU:OE1	3:C:2017:HOH:O	2.02	0.78
1:D:132:VAL:H	1:D:191:ASN:HD22	1.31	0.76
1:E:228:GLN:NE2	1:E:231:ARG:HH11	1.82	0.76
1:E:208[B]:ARG:HG3	1:E:208[B]:ARG:HH11	1.51	0.76
1:E:219:ARG:O	3:E:2156:HOH:O	2.03	0.75
1:F:10:GLU:OE2	3:F:2003:HOH:O	2.08	0.72
1:E:222:ARG:HB3	3:E:2155:HOH:O	1.89	0.71
1:B:59:ARG:NH1	1:B:219:ARG:NH1	2.40	0.70
1:B:32:HIS:CD2	1:B:35:GLY:H	2.11	0.68
1:A:243:HIS:HE1	1:B:223:LYS:NZ	1.93	0.67
1:A:197:GLN:HG3	3:A:2146:HOH:O	1.95	0.65
1:E:89:ASP:OD1	3:E:2081:HOH:O	2.13	0.65
1:E:218:ARG:O	3:E:2155:HOH:O	2.14	0.65
1:E:23:ASP:OD2	1:E:208[A]:ARG:NH2	2.30	0.63
1:E:208[B]:ARG:HG3	1:E:208[B]:ARG:NH1	2.08	0.63
1:E:208[A]:ARG:HH11	1:E:208[A]:ARG:CG	2.13	0.62
1:E:23:ASP:O	1:E:208[B]:ARG:NH2	2.33	0.61
1:D:132:VAL:H	1:D:191:ASN:ND2	1.98	0.61
1:D:6:THR:N	3:D:2001:HOH:O	2.35	0.59
1:B:59:ARG:HH21	1:E:95:PHE:CB	2.16	0.58
1:E:208[B]:ARG:CG	1:E:208[B]:ARG:HH11	2.15	0.58
1:E:208[A]:ARG:HH11	1:E:208[A]:ARG:HG3	1.67	0.58
1:A:243:HIS:HE1	1:B:223:LYS:HZ3	1.50	0.58
1:A:142:ASP:OD2	1:A:145:HIS:HD2	1.87	0.57
1:B:59:ARG:HH11	1:B:219:ARG:NH1	2.01	0.57
1:C:142:ASP:OD2	1:C:145:HIS:HD2	1.88	0.57
1:A:243:HIS:CE1	1:B:223:LYS:HZ3	2.23	0.55
1:E:228:GLN:HE21	1:E:231:ARG:HH11	1.52	0.55
1:F:122:HIS:ND1	1:F:124:GLU:OE2	2.29	0.55
3:C:2169:HOH:O	1:D:86:THR:HB	2.05	0.55
1:B:83:ASN:O	1:B:89:ASP:HB3	2.07	0.54
1:A:243:HIS:CE1	1:B:223:LYS:NZ	2.75	0.54
1:B:239:LEU:O	1:B:243:HIS:HD2	1.91	0.54
1:B:59:ARG:HH11	1:B:219:ARG:CZ	2.21	0.54
1:D:45:HIS:HE1	3:D:2061:HOH:O	1.91	0.54
1:B:196:GLU:OE1	3:B:2139:HOH:O	2.19	0.54
1:B:251:LEU:HD11	3:B:2089:HOH:O	2.09	0.52
1:A:13:GLN:NE2	3:A:2009:HOH:O	2.43	0.51
1:B:167:ASN:HD22	1:B:167:ASN:N	2.10	0.50
1:B:250:ASP:C	1:B:252:GLY:H	2.16	0.49
1:C:157:HIS:HE1	1:C:236:ASP:O	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:223:LYS:HB2	3:E:2156:HOH:O	2.12	0.49
1:E:222:ARG:CB	3:E:2155:HOH:O	2.55	0.49
1:F:221:ALA:O	1:F:225:LEU:HG	2.13	0.49
1:A:250:ASP:HB2	1:E:87:PRO:HB2	1.94	0.49
1:A:7:PRO:HA	3:A:2001:HOH:O	2.11	0.48
1:B:93:ILE:HG23	3:B:2031:HOH:O	2.12	0.48
1:E:228:GLN:NE2	1:E:231:ARG:HD2	2.28	0.48
1:F:14:LYS:NZ	3:F:2008:HOH:O	2.43	0.48
1:C:167:ASN:N	1:C:167:ASN:HD22	2.10	0.48
1:B:136:GLU:HG3	3:B:2078:HOH:O	2.14	0.47
1:B:243:HIS:HE1	1:C:223:LYS:NZ	2.12	0.47
1:E:214:PRO:CG	1:F:251:LEU:HD21	2.44	0.47
1:E:208[B]:ARG:CG	1:E:208[B]:ARG:NH1	2.75	0.46
1:E:41:THR:HG22	1:E:78:ASP:HB3	1.99	0.45
1:B:218:ARG:NH2	3:B:2155:HOH:O	2.43	0.45
1:C:19:ARG:NH1	3:C:2015:HOH:O	2.50	0.45
1:B:251:LEU:HD21	1:C:214:PRO:HG2	1.98	0.44
1:C:8:PHE:HA	1:C:11:TYR:CE2	2.52	0.44
1:A:182:ARG:HD2	3:A:2126:HOH:O	2.17	0.44
1:E:162:HIS:HE1	3:F:2099:HOH:O	2.00	0.44
1:C:82:PHE:HB3	1:C:84:LEU:HD21	2.00	0.44
1:D:37:SER:HB3	1:D:72[A]:SER:OG	2.18	0.44
1:D:142:ASP:OD2	1:D:145:HIS:HD2	2.00	0.44
1:C:162:HIS:HD2	3:C:2041:HOH:O	2.01	0.43
1:B:250:ASP:C	1:B:252:GLY:N	2.72	0.43
1:D:239:LEU:HD23	1:D:239:LEU:C	2.37	0.43
1:C:142:ASP:OD2	1:C:145:HIS:CD2	2.71	0.43
1:B:122:HIS:N	1:B:123:PRO:CD	2.81	0.43
1:B:94:ILE:HG12	1:B:241:LEU:HB3	2.01	0.43
1:B:167:ASN:ND2	1:B:170:ARG:HH21	2.17	0.43
1:E:14:LYS:HE3	1:E:15:TYR:OH	2.19	0.42
1:D:20:LEU:HD23	1:D:29:VAL:HB	2.01	0.42
1:B:218:ARG:NH1	3:B:2155:HOH:O	2.33	0.42
1:E:83:ASN:O	1:E:89:ASP:HB3	2.19	0.42
1:B:145:HIS:HE1	3:B:2171:HOH:O	2.03	0.42
1:B:251:LEU:HD21	1:C:214:PRO:CG	2.50	0.42
1:E:214:PRO:HG2	1:F:251:LEU:HD21	2.02	0.41
1:D:43:THR:O	1:D:47:GLU:HG3	2.20	0.41
1:E:14:LYS:HE2	1:E:54:ASP:OD2	2.20	0.41
1:B:204:TRP:O	1:B:208:ARG:HG2	2.20	0.41
1:C:167:ASN:ND2	1:C:170:ARG:HH21	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:LYS:NZ	3:A:2042:HOH:O	2.53	0.41
1:C:200:LEU:HA	1:C:200:LEU:HD23	1.92	0.41
1:C:221:ALA:O	1:C:225:LEU:HG	2.21	0.40
1:E:200:LEU:N	1:E:201:PRO:CD	2.84	0.40
1:A:143:GLY:N	1:A:144:PRO:CD	2.85	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)
3:B:2142:HOH:O	3:D:2101:HOH:O[1_455]	2.12	0.08

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	247/257 (96%)	244 (99%)	3 (1%)	0	100	100
1	B	247/257 (96%)	244 (99%)	3 (1%)	0	100	100
1	C	246/257 (96%)	241 (98%)	5 (2%)	0	100	100
1	D	248/257 (96%)	244 (98%)	4 (2%)	0	100	100
1	E	248/257 (96%)	246 (99%)	2 (1%)	0	100	100
1	F	246/257 (96%)	242 (98%)	4 (2%)	0	100	100
All	All	1482/1542 (96%)	1461 (99%)	21 (1%)	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	199/215 (93%)	199 (100%)	0	100	100
1	B	202/215 (94%)	199 (98%)	3 (2%)	65	69
1	C	203/215 (94%)	199 (98%)	4 (2%)	55	58
1	D	199/215 (93%)	193 (97%)	6 (3%)	41	41
1	E	204/215 (95%)	198 (97%)	6 (3%)	42	43
1	F	198/215 (92%)	197 (100%)	1 (0%)	88	92
All	All	1205/1290 (93%)	1185 (98%)	20 (2%)	65	65

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	21	GLU
1	B	39	VAL
1	B	216	LEU
1	C	80	THR
1	C	92	GLU
1	C	99	ARG
1	C	167	ASN
1	D	6	THR
1	D	16	GLU
1	D	29	VAL
1	D	63	VAL
1	D	218[A]	ARG
1	D	218[B]	ARG
1	E	4	LEU
1	E	20	LEU
1	E	197	GLN
1	E	208[A]	ARG
1	E	208[B]	ARG
1	E	218	ARG
1	F	213	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	145	HIS
1	A	243	HIS
1	B	32	HIS
1	B	145	HIS
1	B	167	ASN
1	B	243	HIS
1	C	83	ASN
1	C	88	HIS
1	C	102	ASN
1	C	145	HIS
1	C	157	HIS
1	C	162	HIS
1	C	167	ASN
1	D	45	HIS
1	D	145	HIS
1	D	177	GLN
1	D	191	ASN
1	E	162	HIS
1	E	228	GLN
1	F	102	ASN
1	F	103	ASN

5.3.3 RNA

There are no RNA molecules 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 2 ligands modelled in this entry, 2 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

There are no such residues in this entry.

5.8 Polymer linkage issues

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	249:ILE	C	250:ASP	N	1.20

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.