



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2023 – 01:51 AM EDT

PDB ID : 3A2N
Title : Crystal structure of DBJA (Wild Type Type II P21)
Authors : Sato, Y.; Senda, T.
Deposited on : 2009-05-23
Resolution : 1.89 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

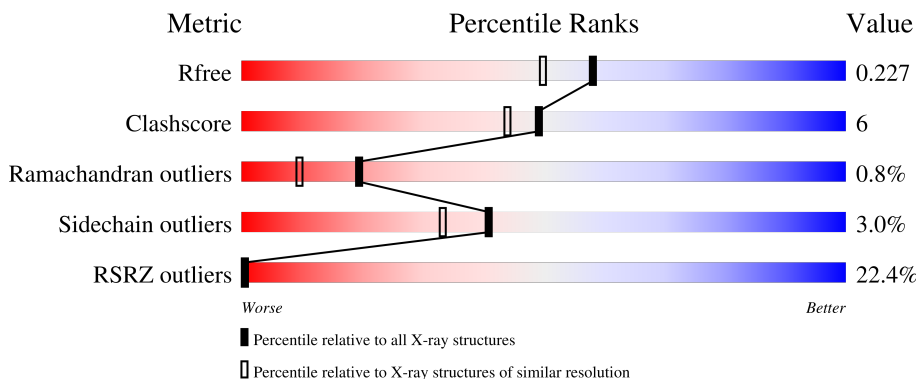
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 $\leq 5\%$. 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	312	 17% 83% 12% . .
1	B	312	 9% 86% 8% . .
1	E	312	 43% 78% 16% . .
1	F	312	 16% 85% 11% . .

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9801 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 Haloalkane dehalogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	298	2320	1485	415	413	7	0	0	0
1	B	298	2330	1491	418	414	7	0	1	0
1	E	298	2320	1485	415	413	7	0	0	0
1	F	299	2338	1497	419	415	7	0	1	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	311	VAL	-	SEE REMARK 999	UNP P59337
A	312	ASP	-	SEE REMARK 999	UNP P59337
B	311	VAL	-	SEE REMARK 999	UNP P59337
B	312	ASP	-	SEE REMARK 999	UNP P59337
E	311	VAL	-	SEE REMARK 999	UNP P59337
E	312	ASP	-	SEE REMARK 999	UNP P59337
F	311	VAL	-	SEE REMARK 999	UNP P59337
F	312	ASP	-	SEE REMARK 999	UNP P59337

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Cl 1	0	0
2	B	1	Total 1	Cl 1	0	0
2	F	1	Total 1	Cl 1	0	0

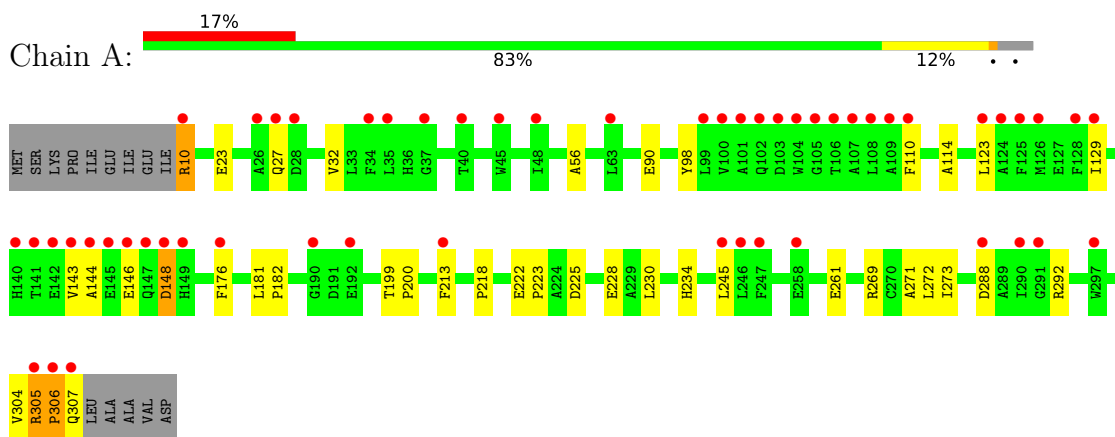
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	110	Total 110	O 110	0	0
3	B	162	Total 162	O 162	0	0
3	E	45	Total 45	O 45	0	0
3	F	172	Total 173	O 173	0	1

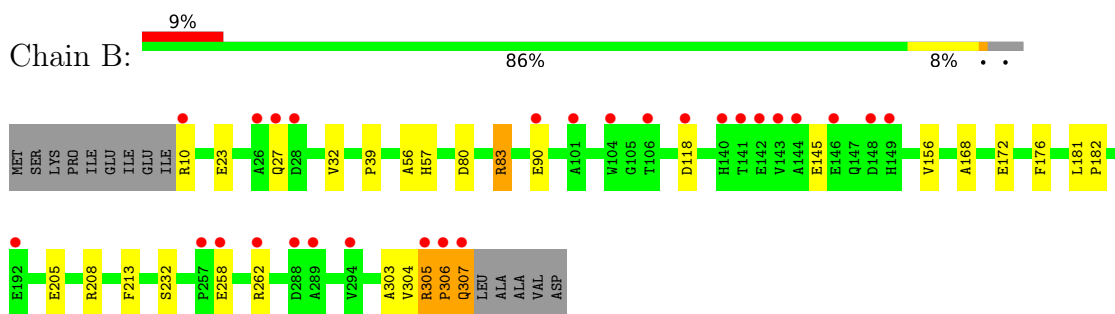
3 Residue-property plots [i](#)

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 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.

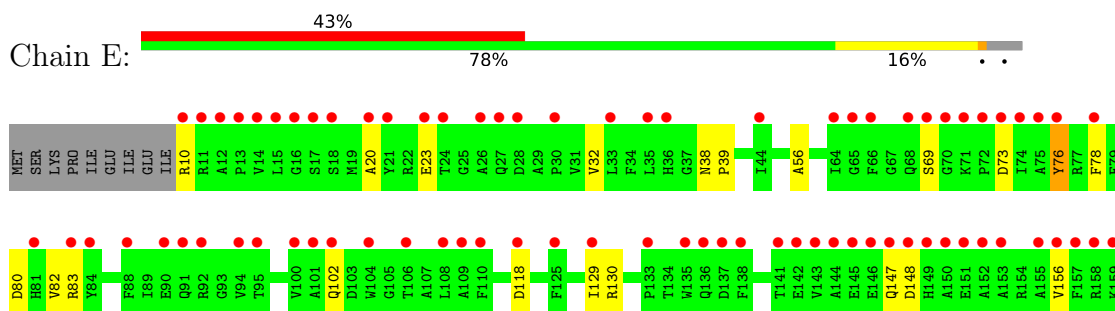
- Molecule 1: Haloalkane dehalogenase

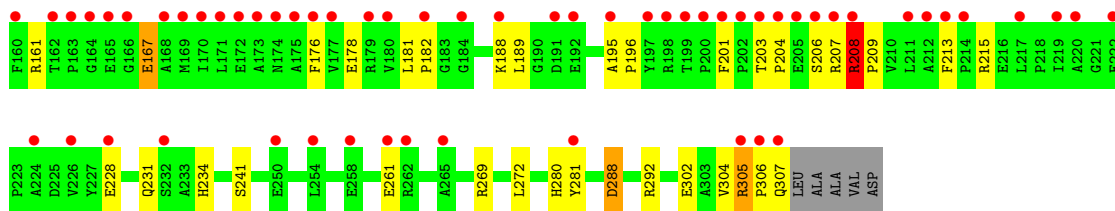


- Molecule 1: Haloalkane dehalogenase

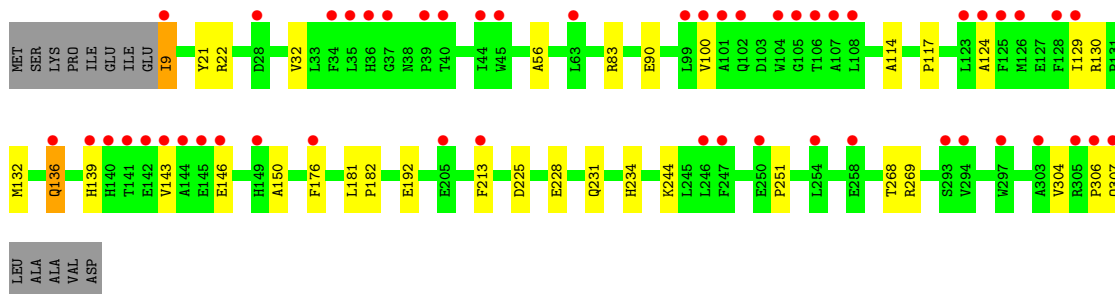
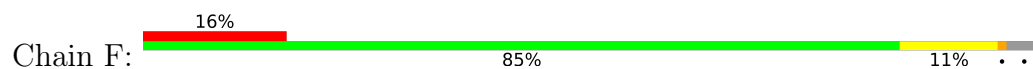


- Molecule 1: Haloalkane dehalogenase





● Molecule 1: Haloalkane dehalogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	125.53Å 47.77Å 99.41Å 90.00° 93.61° 90.00°	Depositor
Resolution (Å)	32.81 – 1.89 31.49 – 1.89	Depositor EDS
% Data completeness (in resolution range)	99.3 (32.81-1.89) 99.3 (31.49-1.89)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.17 (at 1.89Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.192 , 0.228 0.191 , 0.227	Depositor DCC
R_{free} test set	4719 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtrriage
Anisotropy	0.082	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 53.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9801	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.79 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.0884e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: 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	0.59	0/2388	0.61	1/3252 (0.0%)
1	B	0.65	0/2399	0.66	1/3267 (0.0%)
1	E	0.77	6/2388 (0.3%)	0.69	5/3252 (0.2%)
1	F	0.68	0/2407	0.69	2/3278 (0.1%)
All	All	0.68	6/9582 (0.1%)	0.66	9/13049 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	73	ASP	CG-OD2	12.28	1.53	1.25
1	E	76	TYR	C-O	12.03	1.46	1.23
1	E	73	ASP	CG-OD1	10.18	1.48	1.25
1	E	167	GLU	CD-OE1	8.03	1.34	1.25
1	E	208	ARG	CZ-NH1	6.38	1.41	1.33
1	E	76	TYR	CG-CD1	5.92	1.46	1.39

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	73	ASP	CB-CG-OD1	-11.52	107.94	118.30
1	E	76	TYR	CB-CG-CD2	-8.35	115.99	121.00
1	E	80	ASP	CB-CG-OD1	7.92	125.42	118.30
1	E	208	ARG	NE-CZ-NH2	-6.67	116.97	120.30
1	E	76	TYR	CG-CD2-CE2	-6.24	116.31	121.30
1	F	83	ARG	NE-CZ-NH2	-5.91	117.34	120.30
1	A	225	ASP	CB-CG-OD1	5.54	123.29	118.30
1	B	83	ARG	NE-CZ-NH2	-5.25	117.68	120.30
1	F	225	ASP	CB-CG-OD1	5.10	122.89	118.30

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	2320	0	2268	28	0
1	B	2330	0	2274	32	0
1	E	2320	0	2268	38	0
1	F	2338	0	2285	31	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	F	1	0	0	0	0
3	A	110	0	0	1	0
3	B	162	0	0	3	0
3	E	45	0	0	3	0
3	F	173	0	0	2	0
All	All	9801	0	9095	109	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 (109) 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:E:188:LYS:HD3	3:E:327:HOH:O	1.57	1.00
1:B:305:ARG:HB2	1:B:306:PRO:HD2	1.44	1.00
1:E:304:VAL:HG21	1:F:304:VAL:HG23	1.63	0.81
1:F:9:ILE:HG23	1:F:21:TYR:O	1.80	0.81
1:B:307:GLN:HG3	1:F:117:PRO:HB3	1.65	0.79
1:E:304:VAL:CG2	1:F:304:VAL:HG23	2.14	0.78
1:B:83:ARG:HD3	3:B:407:HOH:O	1.85	0.76
1:A:269:ARG:HH22	1:B:307:GLN:NE2	1.87	0.73
1:A:304:VAL:CG2	1:B:304:VAL:HG23	2.20	0.72
1:B:258:GLU:HG2	1:B:262:ARG:NH1	2.04	0.71
1:A:304:VAL:HG21	1:B:304:VAL:HG23	1.73	0.70
1:B:208:ARG:HB2	3:B:396:HOH:O	1.92	0.70
1:F:9:ILE:HG21	1:F:22:ARG:HG3	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:ARG:CB	1:B:306:PRO:HD2	2.22	0.69
1:F:90:GLU:HG2	3:F:433:HOH:O	1.93	0.68
1:B:306:PRO:HA	1:F:114:ALA:O	1.94	0.67
1:A:181:LEU:HB3	1:A:182:PRO:HD3	1.77	0.66
1:B:10:ARG:CZ	1:B:23:GLU:OE1	2.44	0.66
1:E:302:GLU:O	1:E:305:ARG:HG2	1.95	0.66
1:A:304:VAL:CG2	1:B:304:VAL:CG2	2.75	0.64
1:E:304:VAL:O	1:E:305:ARG:O	2.17	0.63
1:E:305:ARG:HB2	1:E:306:PRO:CD	2.29	0.62
1:B:258:GLU:HG2	1:B:262:ARG:HH11	1.64	0.62
1:F:146:GLU:HG3	1:F:251:PRO:CB	2.31	0.60
1:B:181:LEU:HB3	1:B:182:PRO:HD3	1.82	0.60
1:A:269:ARG:NH2	1:B:307:GLN:HB2	2.16	0.60
1:E:188:LYS:HE3	1:E:189:LEU:O	2.01	0.60
1:E:83:ARG:HD2	3:E:350:HOH:O	2.02	0.58
1:F:9:ILE:HG21	1:F:22:ARG:CG	2.33	0.58
1:A:269:ARG:HD3	1:B:303:ALA:O	2.03	0.57
1:E:228:GLU:HA	1:E:231:GLN:HE21	1.68	0.57
1:B:90:GLU:HG2	3:B:362:HOH:O	2.05	0.57
1:E:201:PHE:HB3	1:E:207:ARG:HG2	1.87	0.57
1:E:304:VAL:CG2	1:F:304:VAL:CG2	2.81	0.56
1:B:307:GLN:HG3	1:F:117:PRO:CB	2.35	0.56
1:A:288:ASP:O	1:A:292:ARG:HG2	2.07	0.55
1:E:161:ARG:HA	1:E:215:ARG:HG2	1.89	0.55
1:F:139[B]:HIS:ND1	1:F:150:ALA:HA	2.21	0.55
1:B:80:ASP:OD1	1:B:83:ARG:NH2	2.39	0.54
1:A:304:VAL:HG23	1:B:304:VAL:CG2	2.38	0.54
1:A:114:ALA:O	1:E:307:GLN:N	2.41	0.54
1:F:130:ARG:NH1	3:F:414:HOH:O	2.36	0.54
1:B:32:VAL:HG23	1:B:56:ALA:HB1	1.91	0.53
1:F:268:THR:HG22	1:F:269:ARG:HD2	1.90	0.53
1:B:304:VAL:O	1:B:305:ARG:O	2.26	0.53
1:B:305:ARG:HB2	1:B:306:PRO:CD	2.29	0.53
1:A:222:GLU:HA	1:A:223:PRO:C	2.29	0.53
1:B:307:GLN:N	1:F:114:ALA:O	2.36	0.52
1:B:305:ARG:CB	1:B:306:PRO:CD	2.88	0.52
1:E:102:GLN:NE2	1:E:281:TYR:HA	2.26	0.51
1:E:305:ARG:HB2	1:E:306:PRO:HD2	1.93	0.50
1:A:218:PRO:HG3	1:A:230:LEU:HD12	1.92	0.50
1:E:203:THR:O	1:E:206:SER:N	2.42	0.50
1:E:288:ASP:O	1:E:292:ARG:HG2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:ARG:HB2	1:A:306:PRO:HD2	1.92	0.49
1:E:38:ASN:OD1	1:E:39:PRO:HA	2.11	0.49
1:F:9:ILE:CG2	1:F:22:ARG:HG3	2.42	0.49
1:A:269:ARG:HH22	1:B:307:GLN:HE21	1.57	0.49
1:F:143:VAL:HG11	1:F:146:GLU:HG2	1.94	0.49
1:A:261:GLU:HA	1:A:272:LEU:HD22	1.95	0.49
1:A:90:GLU:HG2	3:A:332:HOH:O	2.13	0.49
1:A:306:PRO:O	1:A:307:GLN:HB2	2.13	0.48
1:F:146:GLU:HG3	1:F:251:PRO:HB3	1.95	0.48
1:B:27:GLN:HG2	1:B:57:HIS:CE1	2.49	0.48
1:A:110:PHE:HE1	1:A:123:LEU:HD21	1.79	0.48
1:F:32:VAL:HG23	1:F:56:ALA:HB1	1.95	0.48
1:A:129:ILE:O	1:A:234:HIS:HE1	1.97	0.47
1:A:148:ASP:N	1:A:148:ASP:OD1	2.48	0.47
1:B:10:ARG:NH1	1:B:23:GLU:OE1	2.48	0.47
1:F:143:VAL:CG1	1:F:146:GLU:HG2	2.45	0.47
1:A:32:VAL:HG22	1:A:98:TYR:HB2	1.97	0.46
1:F:244:LYS:HE2	1:F:268:THR:O	2.15	0.46
1:E:130:ARG:HA	1:E:234:HIS:CE1	2.51	0.46
1:E:102:GLN:NE2	1:E:280:HIS:O	2.48	0.45
1:F:136:GLN:H	1:F:136:GLN:CD	2.21	0.45
1:E:304:VAL:HG23	1:F:304:VAL:CG2	2.47	0.45
1:E:188:LYS:HD2	1:E:189:LEU:N	2.32	0.45
1:E:76:TYR:CE2	1:E:209:PRO:HG3	2.52	0.45
1:E:181:LEU:HB3	1:E:182:PRO:HD3	1.99	0.44
1:E:208:ARG:HB3	1:E:209:PRO:HD3	2.00	0.43
1:A:304:VAL:CG2	1:B:304:VAL:HG21	2.47	0.43
1:F:268:THR:HG22	1:F:269:ARG:CD	2.48	0.43
1:B:168:ALA:O	1:B:172:GLU:HB3	2.18	0.43
1:A:245:LEU:HD12	1:A:271:ALA:HB3	2.01	0.43
1:F:130:ARG:O	1:F:132:MET:HG3	2.19	0.43
1:A:32:VAL:HG23	1:A:56:ALA:HB1	2.01	0.43
1:A:304:VAL:HG23	1:B:304:VAL:HG21	2.00	0.43
1:E:32:VAL:HG23	1:E:56:ALA:HB1	2.01	0.43
1:F:129:ILE:O	1:F:234:HIS:HE1	2.02	0.43
1:E:261:GLU:HA	1:E:272:LEU:HD22	1.99	0.42
1:F:228:GLU:HA	1:F:231:GLN:HE21	1.82	0.42
1:E:78:PHE:O	1:E:82:VAL:HG23	2.19	0.42
1:E:167:GLU:CD	1:E:215:ARG:HH22	2.23	0.42
1:E:195:ALA:HB3	1:E:196:PRO:HD3	2.02	0.42
1:F:181:LEU:HB3	1:F:182:PRO:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:20:ALA:H	1:E:69:SER:HA	1.85	0.42
1:E:10:ARG:CZ	1:E:23:GLU:HB3	2.50	0.42
1:E:102:GLN:HE22	1:E:281:TYR:HA	1.83	0.41
1:B:305:ARG:HG3	1:B:306:PRO:N	2.35	0.41
1:E:129:ILE:O	1:E:234:HIS:HE1	2.03	0.41
1:E:269:ARG:HH21	1:F:307:GLN:HG2	1.85	0.41
1:F:146:GLU:HG3	1:F:251:PRO:HB2	2.03	0.41
1:E:178:GLU:O	1:E:182:PRO:HG2	2.20	0.41
1:F:100:VAL:HA	1:F:124:ALA:O	2.21	0.41
1:A:199:THR:N	1:A:200:PRO:CD	2.84	0.41
1:A:245:LEU:HG	1:A:273:ILE:CD1	2.50	0.41
1:E:118:ASP:HB2	3:E:316:HOH:O	2.20	0.40
1:A:10:ARG:NH1	1:A:23:GLU:H	2.19	0.40
1:E:241:SER:O	1:E:269:ARG:HD2	2.20	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	296/312 (95%)	282 (95%)	12 (4%)	2 (1%)	22	12
1	B	297/312 (95%)	287 (97%)	8 (3%)	2 (1%)	22	12
1	E	296/312 (95%)	281 (95%)	11 (4%)	4 (1%)	11	3
1	F	298/312 (96%)	289 (97%)	8 (3%)	1 (0%)	41	31
All	All	1187/1248 (95%)	1139 (96%)	39 (3%)	9 (1%)	19	9

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	144	ALA

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Mol	Chain	Res	Type
1	B	305	ARG
1	E	148	ASP
1	E	305	ARG
1	F	306	PRO
1	A	306	PRO
1	B	306	PRO
1	E	147	GLN
1	E	204	PRO

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	232/244 (95%)	223 (96%)	9 (4%)	32	23
1	B	233/244 (96%)	224 (96%)	9 (4%)	32	23
1	E	232/244 (95%)	227 (98%)	5 (2%)	52	47
1	F	234/244 (96%)	229 (98%)	5 (2%)	53	48
All	All	931/976 (95%)	903 (97%)	28 (3%)	41	33

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

Mol	Chain	Res	Type
1	A	10	ARG
1	A	27	GLN
1	A	143	VAL
1	A	146	GLU
1	A	148	ASP
1	A	176	PHE
1	A	213	PHE
1	A	228	GLU
1	A	305	ARG
1	B	39	PRO
1	B	118	ASP
1	B	145	GLU
1	B	156	VAL

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Mol	Chain	Res	Type
1	B	176	PHE
1	B	205	GLU
1	B	213	PHE
1	B	232	SER
1	B	307	GLN
1	E	156	VAL
1	E	176	PHE
1	E	208	ARG
1	E	213	PHE
1	E	288	ASP
1	F	9	ILE
1	F	136	GLN
1	F	176	PHE
1	F	192	GLU
1	F	213	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	HIS
1	A	234	HIS
1	B	231	GLN
1	B	234	HIS
1	B	307	GLN
1	E	111	HIS
1	E	139	HIS
1	E	231	GLN
1	E	234	HIS
1	F	231	GLN
1	F	234	HIS

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 3 ligands modelled in this entry, 3 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	298/312 (95%)	1.06	54 (18%) 1 1	20, 26, 39, 55	0
1	B	298/312 (95%)	0.80	27 (9%) 9 10	18, 23, 33, 51	0
1	E	298/312 (95%)	2.17	135 (45%) 0 0	24, 39, 60, 64	0
1	F	299/312 (95%)	0.92	51 (17%) 1 1	15, 20, 33, 58	0
All	All	1193/1248 (95%)	1.24	267 (22%) 0 0	15, 25, 48, 64	0

All (267) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	307	GLN	11.3
1	A	306	PRO	10.5
1	B	306	PRO	9.4
1	B	307	GLN	9.4
1	F	307	GLN	9.3
1	E	203	THR	8.9
1	E	143	VAL	8.6
1	E	211	LEU	8.5
1	F	306	PRO	8.4
1	A	307	GLN	8.0
1	E	306	PRO	7.9
1	E	144	ALA	7.4
1	E	170	ILE	7.1
1	A	148	ASP	6.9
1	E	148	ASP	6.8
1	E	14	VAL	6.8
1	E	205	GLU	6.7
1	E	168	ALA	6.6
1	A	144	ALA	6.2
1	E	16	GLY	6.0
1	E	73	ASP	6.0

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Mol	Chain	Res	Type	RSRZ
1	E	174	ASN	5.9
1	E	160	PHE	5.9
1	E	142	GLU	5.9
1	E	208	ARG	5.9
1	E	70	GLY	5.7
1	B	28	ASP	5.7
1	E	169	MET	5.5
1	E	201	PHE	5.5
1	E	10	ARG	5.4
1	B	305	ARG	5.3
1	E	206	SER	5.2
1	E	176	PHE	5.2
1	E	173	ALA	5.2
1	E	84	TYR	5.2
1	E	171	LEU	5.2
1	A	145	GLU	5.1
1	A	28	ASP	5.1
1	E	197	TYR	5.1
1	E	74	ILE	5.0
1	E	164	GLY	5.0
1	E	162	THR	5.0
1	A	142	GLU	4.9
1	E	219	ILE	4.9
1	E	200	PRO	4.8
1	A	26	ALA	4.7
1	E	159	LYS	4.7
1	A	100	VAL	4.6
1	B	27	GLN	4.6
1	A	146	GLU	4.6
1	A	35	LEU	4.5
1	F	35	LEU	4.5
1	E	217	LEU	4.4
1	E	76	TYR	4.4
1	A	27	GLN	4.4
1	F	305	ARG	4.4
1	E	65	GLY	4.4
1	E	179	ARG	4.4
1	E	152	ALA	4.3
1	A	143	VAL	4.3
1	E	191	ASP	4.3
1	E	18	SER	4.2
1	E	27	GLN	4.2

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Mol	Chain	Res	Type	RSRZ
1	E	202	PRO	4.2
1	E	28	ASP	4.1
1	E	150	ALA	4.1
1	A	305	ARG	4.1
1	F	144	ALA	4.1
1	E	226	VAL	4.1
1	E	172	GLU	4.1
1	E	145	GLU	4.0
1	E	204	PRO	4.0
1	E	258	GLU	4.0
1	A	147	GLN	4.0
1	B	26	ALA	4.0
1	F	125	PHE	4.0
1	A	125	PHE	3.9
1	B	118	ASP	3.8
1	E	146	GLU	3.8
1	E	17	SER	3.8
1	E	13	PRO	3.8
1	A	104	TRP	3.8
1	E	133	PRO	3.8
1	B	144	ALA	3.8
1	F	101	ALA	3.7
1	F	100	VAL	3.7
1	F	258	GLU	3.7
1	E	184	GLY	3.6
1	E	136	GLN	3.6
1	E	104	TRP	3.6
1	A	37	GLY	3.6
1	A	124	ALA	3.6
1	E	12	ALA	3.5
1	E	213	PHE	3.5
1	A	101	ALA	3.5
1	E	92	ARG	3.5
1	A	149	HIS	3.5
1	F	28	ASP	3.5
1	A	106	THR	3.5
1	E	24	THR	3.5
1	E	157	PHE	3.5
1	E	305	ARG	3.4
1	E	175	ALA	3.4
1	E	21	TYR	3.4
1	E	149	HIS	3.4

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Mol	Chain	Res	Type	RSRZ
1	F	104	TRP	3.4
1	E	15	LEU	3.4
1	B	143	VAL	3.4
1	E	35	LEU	3.3
1	E	199	THR	3.3
1	E	78	PHE	3.3
1	A	245	LEU	3.3
1	E	214	PRO	3.3
1	F	143	VAL	3.3
1	E	220	ALA	3.3
1	B	149	HIS	3.3
1	F	126	MET	3.3
1	F	37	GLY	3.2
1	B	142	GLU	3.2
1	F	146	GLU	3.2
1	E	163	PRO	3.2
1	E	94	VAL	3.2
1	E	109	ALA	3.2
1	A	141	THR	3.1
1	F	63	LEU	3.1
1	B	141	THR	3.1
1	A	109	ALA	3.1
1	F	106	THR	3.1
1	F	254	LEU	3.1
1	B	262	ARG	3.0
1	A	34	PHE	3.0
1	A	63	LEU	3.0
1	F	9	ILE	3.0
1	E	261	GLU	3.0
1	A	126	MET	3.0
1	E	129	ILE	3.0
1	E	147	GLN	3.0
1	B	140	HIS	3.0
1	E	95	THR	3.0
1	F	123	LEU	2.9
1	B	288	ASP	2.9
1	F	129	ILE	2.9
1	F	145	GLU	2.9
1	A	213	PHE	2.9
1	E	212	ALA	2.9
1	A	105	GLY	2.8
1	E	75	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	205	GLU	2.8
1	A	99	LEU	2.8
1	A	258	GLU	2.8
1	A	247	PHE	2.8
1	E	66	PHE	2.8
1	F	108	LEU	2.8
1	E	166	GLY	2.8
1	A	10	ARG	2.8
1	A	192	GLU	2.8
1	E	20	ALA	2.7
1	F	303	ALA	2.7
1	E	108	LEU	2.7
1	E	254	LEU	2.7
1	F	149	HIS	2.7
1	E	141	THR	2.7
1	E	88	PHE	2.7
1	B	90	GLU	2.7
1	E	72	PRO	2.7
1	B	258	GLU	2.7
1	E	265	ALA	2.6
1	A	288	ASP	2.6
1	B	148	ASP	2.6
1	F	247	PHE	2.6
1	E	180	VAL	2.6
1	E	91	GLN	2.6
1	E	118	ASP	2.6
1	E	71	LYS	2.6
1	E	207	ARG	2.6
1	E	138	PHE	2.6
1	E	177	VAL	2.6
1	F	294	VAL	2.6
1	A	107	ALA	2.6
1	E	155	ALA	2.6
1	E	44	ILE	2.6
1	F	136	GLN	2.6
1	F	34	PHE	2.5
1	E	101	ALA	2.5
1	B	257	PRO	2.5
1	F	297	TRP	2.5
1	F	140	HIS	2.5
1	A	128	PHE	2.5
1	E	110	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	E	100	VAL	2.5
1	E	188	LYS	2.5
1	A	40	THR	2.5
1	A	190	GLY	2.5
1	F	105	GLY	2.5
1	F	99	LEU	2.5
1	E	232	SER	2.5
1	A	48	ILE	2.4
1	E	228	GLU	2.4
1	A	176	PHE	2.4
1	A	103	ASP	2.4
1	E	198	ARG	2.4
1	E	262	ARG	2.4
1	F	102	GLN	2.4
1	A	45	TRP	2.4
1	E	23	GLU	2.4
1	A	129	ILE	2.4
1	E	156	VAL	2.4
1	F	39	PRO	2.4
1	F	213	PHE	2.4
1	E	195	ALA	2.4
1	E	224	ALA	2.4
1	A	108	LEU	2.3
1	E	36	HIS	2.3
1	E	11	ARG	2.3
1	E	151	GLU	2.3
1	E	102	GLN	2.3
1	F	107	ALA	2.3
1	F	176	PHE	2.3
1	A	297	TRP	2.3
1	E	81	HIS	2.3
1	F	40	THR	2.3
1	A	110	PHE	2.3
1	F	44	ILE	2.3
1	E	182	PRO	2.3
1	A	246	LEU	2.3
1	E	158	ARG	2.3
1	B	104	TRP	2.2
1	E	68	GLN	2.2
1	E	90	GLU	2.2
1	E	165	GLU	2.2
1	E	83	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	64	ILE	2.2
1	E	69	SER	2.2
1	F	246	LEU	2.2
1	E	30	PRO	2.2
1	E	222	GLU	2.2
1	A	102	GLN	2.2
1	F	128	PHE	2.2
1	B	294	VAL	2.2
1	F	142	GLU	2.2
1	F	124	ALA	2.2
1	B	192	GLU	2.2
1	F	293	SER	2.1
1	F	36	HIS	2.1
1	A	123	LEU	2.1
1	B	10	ARG	2.1
1	B	106	THR	2.1
1	F	139[A]	HIS	2.1
1	E	153	ALA	2.1
1	E	281	TYR	2.1
1	E	106	THR	2.1
1	E	192	GLU	2.1
1	E	137	ASP	2.1
1	E	125	PHE	2.1
1	F	250	GLU	2.1
1	E	33	LEU	2.1
1	A	291	GLY	2.0
1	A	290	ILE	2.0
1	E	250	GLU	2.0
1	F	141	THR	2.0
1	A	140	HIS	2.0
1	B	101	ALA	2.0
1	B	289	ALA	2.0
1	E	26	ALA	2.0
1	E	135	TRP	2.0
1	F	45	TRP	2.0
1	B	146	GLU	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 [i](#)

There are no monosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	CL	A	313	1/1	0.99	0.04	15,15,15,15	0
2	CL	B	313	1/1	0.99	0.05	12,12,12,12	0
2	CL	F	313	1/1	1.00	0.06	10,10,10,10	0

6.5 Other polymers [i](#)

There are no such residues in this entry.