



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 10:24 PM UTC

PDB ID : 4ARX / pdb_00004arx
Title : Lepidoptera-specific toxin Cry1Ac from *Bacillus thuringiensis* ssp. *kurstaki* HD-73
Authors : Derbyshire, D.J.; Carroll, J.; Ellar, D.J.; Li, J.
Deposited on : 2012-04-27
Resolution : 2.35 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

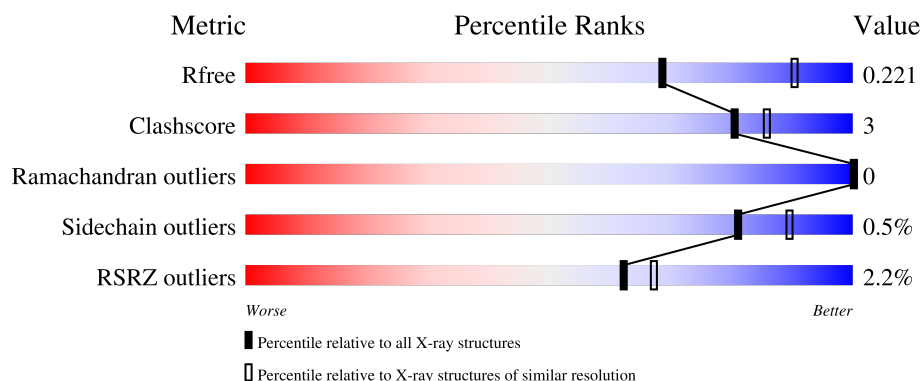
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	180053	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	579	94% 6%
1	B	579	94% 5%
1	C	579	4% 93% 7%
1	D	579	3% 92% 7%

2 Entry composition [i](#)

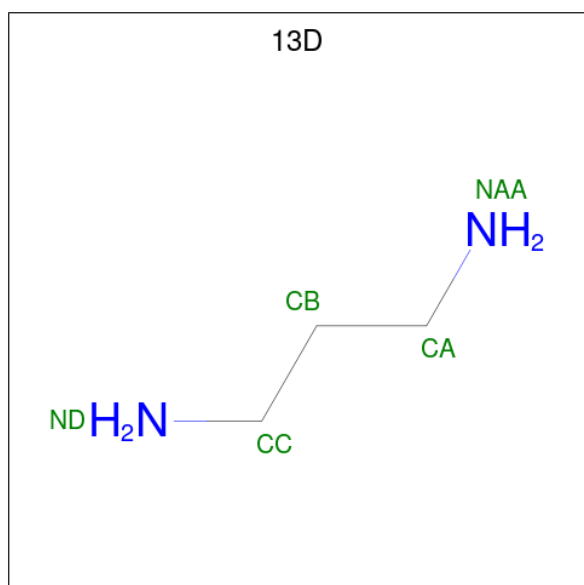
There are 4 unique types of molecules in this entry. The entry contains 19105 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PESTICIDAL CRYSTAL PROTEIN CRY1AC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	579	Total	C	N	O	S	0	0	1
			4594	2924	804	859	7			
1	B	578	Total	C	N	O	S	0	0	0
			4594	2924	803	860	7			
1	C	578	Total	C	N	O	S	0	0	1
			4587	2920	803	857	7			
1	D	577	Total	C	N	O	S	0	0	1
			4583	2918	802	856	7			

- Molecule 2 is 1,3-DIAMINOPROPANE (CCD ID: 13D) (formula: C₃H₁₀N₂).



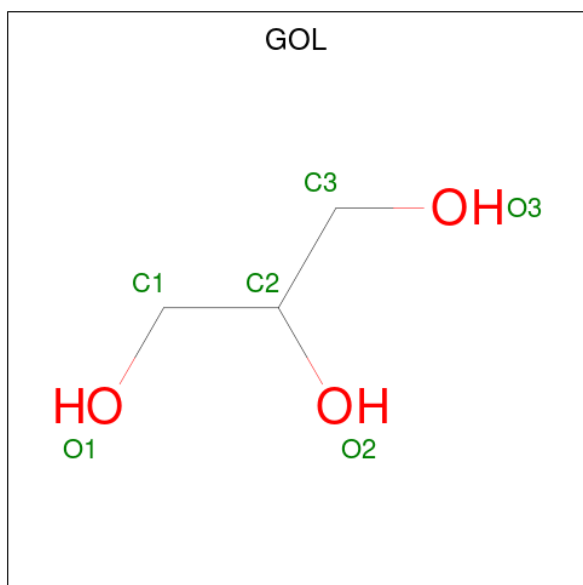
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			5	3	2		
2	A	1	Total	C	N	0	0
			5	3	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	N	0	0
			5	3	2		
2	B	1	Total	C	N	0	0
			5	3	2		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			6	3	3		

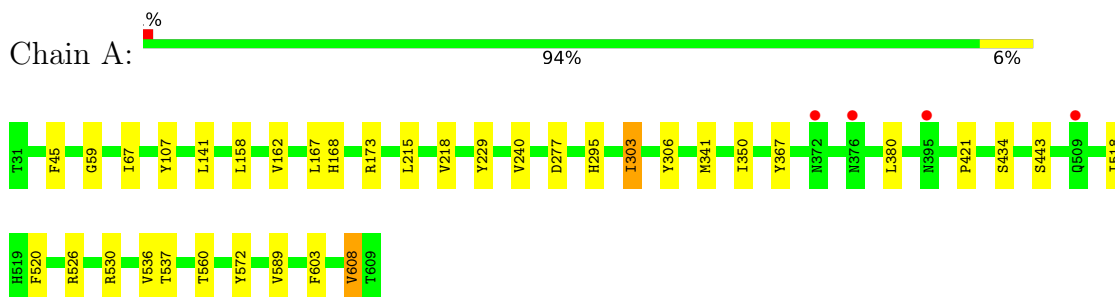
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	189	Total	O	0	0
			189	189		
4	B	208	Total	O	0	0
			208	208		
4	C	163	Total	O	0	0
			163	163		
4	D	161	Total	O	0	0
			161	161		

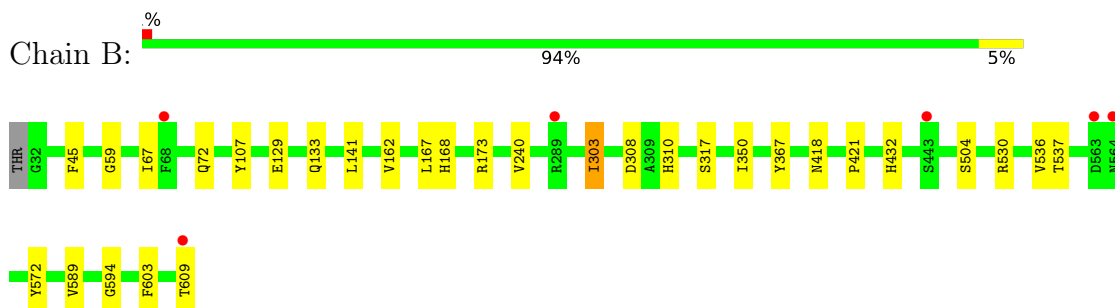
3 Residue-property plots [i](#)

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.

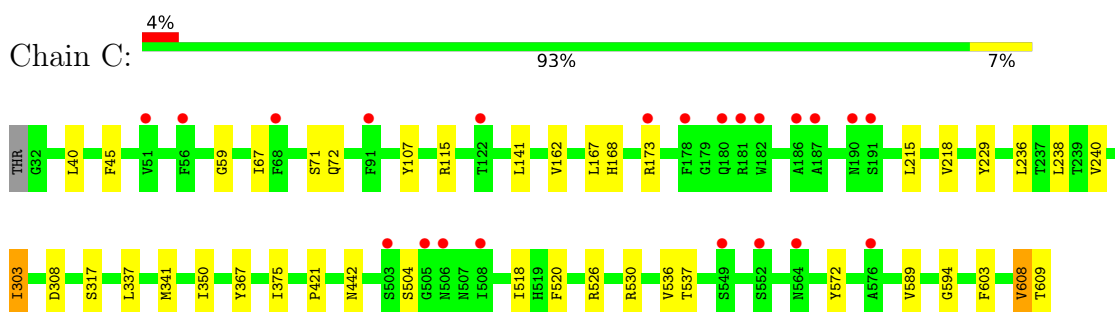
- Molecule 1: PESTICIDAL CRYSTAL PROTEIN CRY1AC



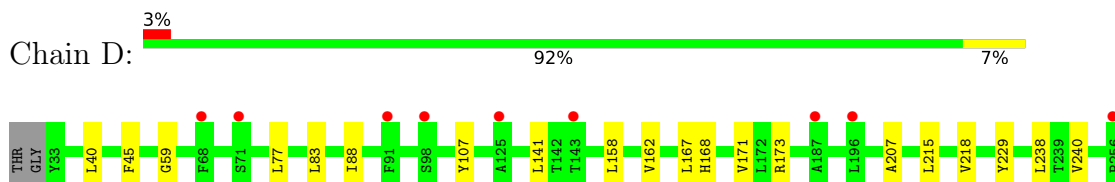
- Molecule 1: PESTICIDAL CRYSTAL PROTEIN CRY1AC

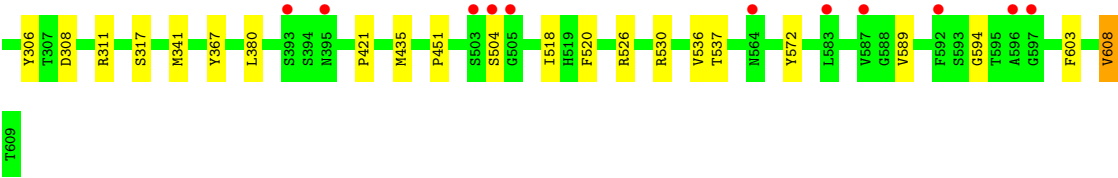


- Molecule 1: PESTICIDAL CRYSTAL PROTEIN CRY1AC



- Molecule 1: PESTICIDAL CRYSTAL PROTEIN CRY1AC





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	51.81Å 112.97Å 123.30Å 113.11° 91.52° 100.51°	Depositor
Resolution (Å)	112.51 – 2.35 112.51 – 2.35	Depositor EDS
% Data completeness (in resolution range)	91.7 (112.51-2.35) 91.7 (112.51-2.35)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.184 , 0.218 0.187 , 0.221	Depositor DCC
R_{free} test set	4815 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	22.3	Xtriage
Anisotropy	0.376	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.035 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	19105	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.77% of the height of the origin peak. No significant pseudotranslation is detected.*

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: 13D, GOL

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	0.76	2/4714 (0.0%)	0.78	0/6425
1	B	0.78	3/4714 (0.1%)	0.80	1/6423 (0.0%)
1	C	0.75	2/4707 (0.0%)	0.79	2/6415 (0.0%)
1	D	0.72	2/4703 (0.0%)	0.77	0/6410
All	All	0.75	9/18838 (0.0%)	0.79	3/25673 (0.0%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	421	PRO	CA-C	8.05	1.56	1.51
1	B	421	PRO	CA-C	7.81	1.56	1.51
1	A	608	VAL	C-N	-7.32	1.23	1.33
1	C	608	VAL	C-N	-7.10	1.23	1.33
1	D	421	PRO	CA-C	6.54	1.55	1.51
1	D	608	VAL	C-N	-6.46	1.24	1.33
1	B	432	HIS	CG-CD2	6.14	1.42	1.35
1	C	421	PRO	CA-C	5.85	1.55	1.51
1	B	310	HIS	CG-ND1	-5.01	1.32	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	71	SER	N-CA-C	5.41	117.60	111.11
1	B	367	TYR	N-CA-C	5.17	117.54	107.71
1	C	367	TYR	N-CA-C	5.00	117.21	107.71

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	4594	0	4448	28	0
1	B	4594	0	4448	17	0
1	C	4587	0	4441	27	1
1	D	4583	0	4438	22	1
2	A	10	0	20	1	0
2	B	10	0	20	0	0
3	D	6	0	8	0	0
4	A	189	0	0	5	0
4	B	208	0	0	1	0
4	C	163	0	0	1	0
4	D	161	0	0	0	0
All	All	19105	0	17823	93	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 3.

All (93) 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:608:VAL:HG23	1:C:609:THR:N	1.96	0.80
1:D:530:ARG:HD3	1:D:572:TYR:CD2	2.24	0.72
1:C:608:VAL:CG2	1:C:609:THR:N	2.54	0.71
1:B:530:ARG:HD3	1:B:572:TYR:CD2	2.26	0.70
1:A:67:ILE:CD1	4:A:2026:HOH:O	2.40	0.69
1:A:530:ARG:HD3	1:A:572:TYR:CD2	2.28	0.68
1:A:162:VAL:HG21	1:A:240:VAL:HG13	1.74	0.68
1:C:530:ARG:HD3	1:C:572:TYR:CD2	2.29	0.67
1:C:107:TYR:CE1	1:C:167:LEU:HD11	2.34	0.63
1:A:536:VAL:HG23	1:A:537:THR:HG23	1.80	0.62
1:B:536:VAL:HG23	1:B:537:THR:HG23	1.80	0.62
1:D:530:ARG:HD3	1:D:572:TYR:HD2	1.64	0.62
1:C:536:VAL:HG22	4:C:2154:HOH:O	2.00	0.61
1:B:107:TYR:CE1	1:B:167:LEU:HD11	2.36	0.60
1:C:536:VAL:HG23	1:C:537:THR:HG23	1.83	0.60
1:B:162:VAL:HG21	1:B:240:VAL:HG13	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:526:ARG:HB2	1:A:608:VAL:HG22	1.82	0.60
1:B:530:ARG:HD3	1:B:572:TYR:HD2	1.66	0.60
1:C:162:VAL:HG21	1:C:240:VAL:HG13	1.84	0.59
1:A:67:ILE:HD11	4:A:2026:HOH:O	2.00	0.59
1:C:530:ARG:HD3	1:C:572:TYR:HD2	1.68	0.59
1:B:418:ASN:ND2	4:B:2002:HOH:O	2.36	0.58
1:D:141:LEU:HD13	1:D:168:HIS:HA	1.86	0.57
1:A:530:ARG:HD3	1:A:572:TYR:HD2	1.69	0.56
1:B:303:ILE:HG23	1:B:350:ILE:HB	1.88	0.55
1:A:67:ILE:HD12	1:A:67:ILE:C	2.32	0.54
1:D:526:ARG:HB2	1:D:608:VAL:HG22	1.88	0.54
1:A:560:THR:HG23	4:A:2180:HOH:O	2.07	0.54
1:C:107:TYR:CD1	1:C:167:LEU:HD11	2.42	0.54
1:C:337:LEU:HD22	1:C:341:MET:HE3	1.90	0.54
1:B:107:TYR:CD1	1:B:167:LEU:HD11	2.44	0.52
1:C:303:ILE:HG23	1:C:350:ILE:HB	1.90	0.52
1:A:277:ASP:HB2	2:A:1610:13D:HDN1	1.75	0.52
1:A:303:ILE:HG23	1:A:350:ILE:HB	1.91	0.52
1:C:141:LEU:HD13	1:C:168:HIS:HA	1.91	0.52
1:A:443:SER:OG	1:C:115:ARG:NH1	2.42	0.52
1:A:67:ILE:HD12	1:A:67:ILE:O	2.10	0.51
1:A:141:LEU:HD13	1:A:168:HIS:HA	1.93	0.51
1:C:67:ILE:HD12	1:C:72:GLN:HB3	1.92	0.50
1:C:526:ARG:HB2	1:C:608:VAL:HG22	1.92	0.50
1:A:306:TYR:CZ	1:A:341:MET:HE2	2.46	0.50
1:D:518:ILE:HD12	1:D:520:PHE:CZ	2.46	0.50
1:C:67:ILE:CD1	1:C:72:GLN:HB3	2.43	0.49
1:A:518:ILE:HD12	1:A:520:PHE:CZ	2.47	0.49
1:B:67:ILE:HD12	1:B:72:GLN:HB3	1.94	0.49
1:C:67:ILE:HD13	1:C:72:GLN:NE2	2.28	0.49
1:D:107:TYR:CD1	1:D:167:LEU:HD11	2.47	0.48
1:B:67:ILE:HD13	1:B:72:GLN:NE2	2.29	0.48
1:A:526:ARG:HB2	1:A:608:VAL:CG2	2.42	0.47
1:D:589:VAL:HG23	1:D:603:PHE:CE2	2.50	0.47
1:C:40:LEU:HD21	1:C:238:LEU:HD22	1.97	0.46
1:C:518:ILE:HD12	1:C:520:PHE:CZ	2.51	0.46
1:D:77:LEU:HD13	1:D:88:ILE:HD12	1.98	0.46
1:D:45:PHE:CZ	1:D:59:GLY:HA3	2.51	0.46
1:A:218:VAL:HG12	1:A:218:VAL:O	2.15	0.46
1:B:589:VAL:HG23	1:B:603:PHE:CE2	2.51	0.45
1:D:107:TYR:CE1	1:D:167:LEU:HD11	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:589:VAL:HG23	1:A:603:PHE:CE2	2.51	0.45
1:A:215:LEU:HG	1:A:229:TYR:CE1	2.52	0.44
1:A:295:HIS:HE1	4:A:2086:HOH:O	1.99	0.44
1:A:107:TYR:CE1	1:A:167:LEU:HD11	2.52	0.44
1:C:67:ILE:HD13	1:C:72:GLN:CD	2.42	0.44
1:C:218:VAL:O	1:C:218:VAL:HG12	2.17	0.44
1:C:215:LEU:HG	1:C:229:TYR:CE1	2.53	0.44
1:A:434:SER:HB2	4:A:2094:HOH:O	2.16	0.44
1:D:306:TYR:CZ	1:D:341:MET:HE2	2.53	0.44
1:A:218:VAL:O	1:A:218:VAL:CG1	2.66	0.44
1:B:141:LEU:HD13	1:B:168:HIS:HA	1.99	0.44
1:D:162:VAL:HG21	1:D:240:VAL:HG13	2.00	0.43
1:C:308:ASP:HB2	1:C:317:SER:HB3	2.01	0.43
1:A:45:PHE:CZ	1:A:59:GLY:HA3	2.54	0.43
1:D:40:LEU:HD21	1:D:238:LEU:HD22	2.01	0.43
1:B:67:ILE:HD13	1:B:72:GLN:CD	2.45	0.42
1:B:308:ASP:HB2	1:B:317:SER:HB3	2.02	0.42
1:D:589:VAL:HG23	1:D:603:PHE:CZ	2.55	0.42
1:C:589:VAL:HG23	1:C:603:PHE:CE2	2.54	0.42
1:B:45:PHE:CZ	1:B:59:GLY:HA3	2.54	0.42
1:D:308:ASP:HB2	1:D:317:SER:HB3	2.02	0.42
1:D:215:LEU:HG	1:D:229:TYR:CE1	2.55	0.42
1:D:536:VAL:HG23	1:D:537:THR:HG23	2.02	0.42
1:D:218:VAL:O	1:D:218:VAL:HG12	2.19	0.41
1:A:367:TYR:HA	1:A:380:LEU:O	2.21	0.41
1:D:158:LEU:HD11	1:D:207:ALA:HA	2.03	0.41
1:A:589:VAL:HG23	1:A:603:PHE:CZ	2.55	0.41
1:C:375:ILE:HD12	1:C:375:ILE:HG23	1.76	0.41
1:C:504:SER:OG	1:C:594:GLY:HA2	2.21	0.41
1:D:367:TYR:HA	1:D:380:LEU:O	2.20	0.41
1:D:435:MET:HE2	1:D:451:PRO:HG3	2.02	0.41
1:B:129:GLU:O	1:B:133:GLN:HG2	2.21	0.40
1:A:158:LEU:O	1:A:162:VAL:HG23	2.22	0.40
1:D:504:SER:OG	1:D:594:GLY:HA2	2.22	0.40
1:C:45:PHE:CZ	1:C:59:GLY:HA3	2.55	0.40
1:B:504:SER:OG	1:B:594:GLY:HA2	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)
1:C:442:ASN:OD1	1:D:311:ARG:NH2[1_556]	2.19	0.01

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	577/579 (100%)	564 (98%)	13 (2%)	0	100	100
1	B	576/579 (100%)	561 (97%)	15 (3%)	0	100	100
1	C	576/579 (100%)	561 (97%)	15 (3%)	0	100	100
1	D	575/579 (99%)	559 (97%)	16 (3%)	0	100	100
All	All	2304/2316 (100%)	2245 (97%)	59 (3%)	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	501/502 (100%)	499 (100%)	2 (0%)	84	91
1	B	501/502 (100%)	498 (99%)	3 (1%)	78	87
1	C	500/502 (100%)	497 (99%)	3 (1%)	78	87
1	D	500/502 (100%)	497 (99%)	3 (1%)	78	87
All	All	2002/2008 (100%)	1991 (100%)	11 (0%)	81	89

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	173	ARG
1	A	303	ILE
1	B	173	ARG
1	B	303	ILE
1	B	609	THR
1	C	173	ARG
1	C	236	LEU
1	C	303	ILE
1	D	83	LEU
1	D	171	VAL
1	D	173	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	GLN
1	A	356	GLN
1	A	416	GLN
1	A	566	GLN
1	B	79	GLN
1	B	347	GLN
1	B	416	GLN
1	B	418	ASN
1	C	343	ASN
1	C	416	GLN
1	C	566	GLN
1	D	356	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry

5 ligands are modelled in this entry.

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
2	13D	A	1610	-	4,4,4	0.45	0	3,3,3	0.44	0
2	13D	B	1610	-	4,4,4	0.44	0	3,3,3	0.19	0
3	GOL	D	1609	-	5,5,5	0.12	0	5,5,5	0.63	0
2	13D	A	1609	-	4,4,4	0.49	0	3,3,3	0.65	0
2	13D	B	1611	-	4,4,4	0.44	0	3,3,3	0.28	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	13D	A	1610	-	-	0/2/2/2	-
2	13D	B	1610	-	-	2/2/2/2	-
3	GOL	D	1609	-	-	2/4/4/4	-
2	13D	A	1609	-	-	0/2/2/2	-
2	13D	B	1611	-	-	2/2/2/2	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1609	GOL	O1-C1-C2-O2
3	D	1609	GOL	O1-C1-C2-C3
2	B	1610	13D	CA-CB-CC-ND
2	B	1611	13D	CA-CB-CC-ND
2	B	1610	13D	NAA-CA-CB-CC

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Mol	Chain	Res	Type	Atoms
2	B	1611	13D	NAA-CA-CB-CC

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1610	13D	1	0

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	579/579 (100%)	-0.15	4 (0%)	84 87	11, 21, 36, 57	0
1	B	578/579 (99%)	-0.11	6 (1%)	79 83	9, 19, 33, 54	0
1	C	578/579 (99%)	0.23	22 (3%)	44 50	10, 26, 43, 65	0
1	D	577/579 (99%)	0.41	20 (3%)	47 53	13, 29, 47, 62	0
All	All	2312/2316 (99%)	0.09	52 (2%)	62 67	9, 23, 43, 65	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	506	ASN	4.0
1	C	187	ALA	4.0
1	B	609	THR	3.9
1	D	187	ALA	3.2
1	A	372	ASN	3.2
1	C	182	TRP	3.1
1	C	508	ILE	3.0
1	C	68	PHE	3.0
1	B	68	PHE	2.9
1	C	186	ALA	2.9
1	C	505	GLY	2.9
1	C	178	PHE	2.9
1	D	68	PHE	2.9
1	C	564	ASN	2.7
1	D	505	GLY	2.7
1	A	376	ASN	2.7
1	D	592	PHE	2.6
1	C	180	GLN	2.6
1	C	51	VAL	2.6
1	D	91	PHE	2.6
1	C	549	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	289	ARG	2.5
1	B	443	SER	2.5
1	C	191	SER	2.5
1	C	552	SER	2.5
1	D	71	SER	2.4
1	D	395	ASN	2.4
1	D	256	PRO	2.4
1	D	504	SER	2.4
1	D	596	ALA	2.4
1	C	91	PHE	2.3
1	B	563	ASP	2.3
1	C	190	ASN	2.3
1	D	393	SER	2.3
1	B	564	ASN	2.3
1	D	597	GLY	2.3
1	A	509	GLN	2.3
1	C	181	ARG	2.3
1	D	143	THR	2.2
1	D	196	LEU	2.2
1	D	125	ALA	2.2
1	D	587	VAL	2.2
1	A	395	ASN	2.2
1	C	173	ARG	2.2
1	C	122	THR	2.2
1	D	98	SER	2.2
1	D	503	SER	2.2
1	D	564	ASN	2.1
1	C	576	ALA	2.1
1	D	583	LEU	2.1
1	C	56	PHE	2.1
1	C	503	SER	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	GOL	D	1609	6/6	0.75	0.17	47,50,51,55	0
2	13D	B	1610	5/5	0.88	0.12	27,28,29,29	0
2	13D	A	1610	5/5	0.89	0.10	21,21,22,22	0
2	13D	B	1611	5/5	0.90	0.11	23,25,25,26	0
2	13D	A	1609	5/5	0.92	0.10	10,10,10,11	0

6.5 Other polymers [i](#)

There are no such residues in this entry.