



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:00 PM UTC

PDB ID : 3K44 / pdb_00003k44
Title : Crystal Structure of Drosophila melanogaster Pur-alpha
Authors : Graebisch, A.; Roche, S.; Niessing, D.
Deposited on : 2009-10-05
Resolution : 2.10 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

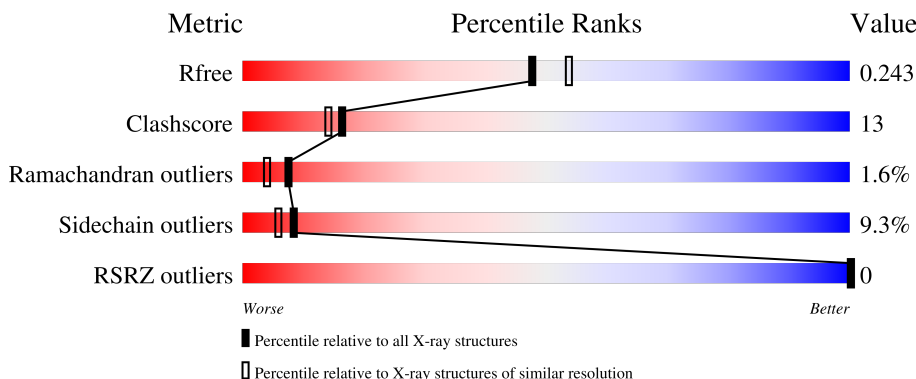
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

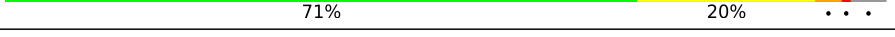
The reported resolution of this entry is 2.10 Å.

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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	146	 75% 16% 5% . .
1	B	146	 72% 19% . . .
1	C	146	 75% 16% 5% . .
1	D	146	 71% 20% . . .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4834 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 Purine-rich binding protein-alpha, isoform B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	145	Total	C	N	O	S	0	1	0
			1162	731	205	222	4			
1	B	140	Total	C	N	O	S	0	4	0
			1156	732	203	217	4			
1	C	145	Total	C	N	O	S	0	2	0
			1174	738	208	224	4			
1	D	140	Total	C	N	O	S	0	5	0
			1162	736	203	219	4			

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0

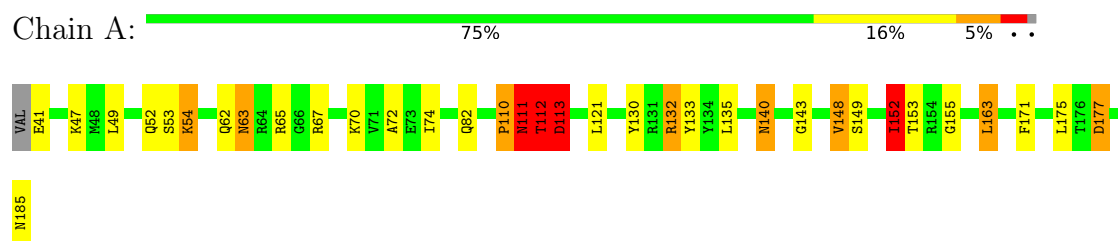
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	45	Total O 45 45	0	0
4	B	39	Total O 39 39	0	0
4	C	42	Total O 42 42	0	0
4	D	30	Total O 30 30	0	0

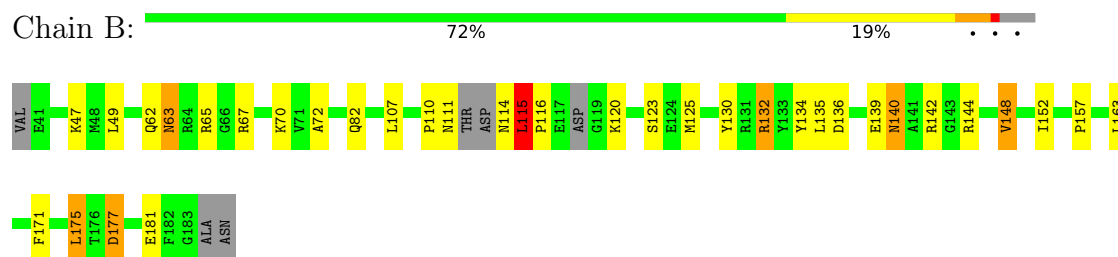
3 Residue-property plots

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.

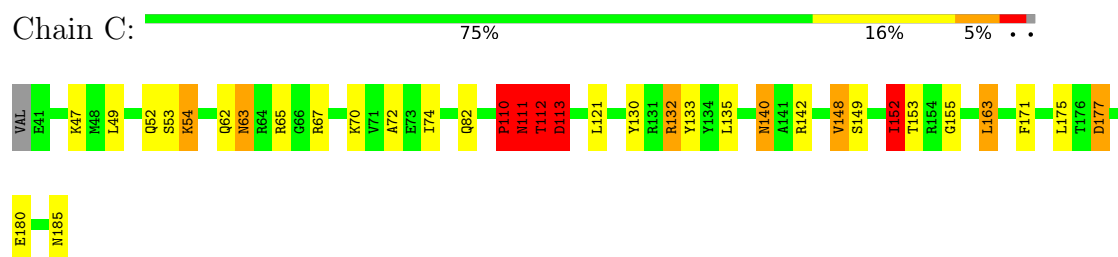
- Molecule 1: Purine-rich binding protein-alpha, isoform B



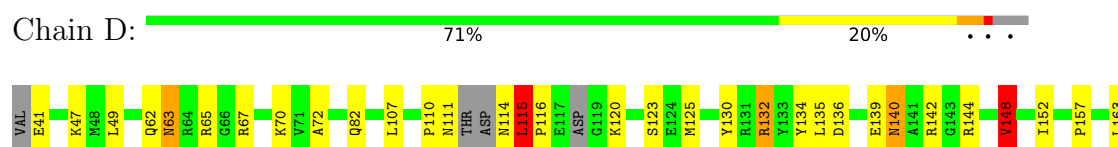
- Molecule 1: Purine-rich binding protein-alpha, isoform B



- Molecule 1: Purine-rich binding protein-alpha, isoform B



- Molecule 1: Purine-rich binding protein-alpha, isoform B



M168	F171	L175	T176	D177	G183	ALA	ASN
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	63.17Å 62.54Å 64.48Å 90.00° 90.01° 90.00°	Depositor
Resolution (Å)	44.46 – 2.10 44.46 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.9 (44.46-2.10) 99.5 (44.46-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.73 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.222 , 0.240 0.226 , 0.243	Depositor DCC
R_{free} test set	1500 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 31.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.027 for l,k,-h 0.026 for -h,-l,-k 0.017 for -h,l,k 0.024 for k,h,-l 0.022 for -k,-h,-l 0.003 for l,h,k 0.004 for k,l,h 0.004 for -l,-h,k 0.004 for -k,-l,h 0.480 for h,-k,-l 0.028 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4834	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.20% 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

5.1 Standard geometry

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

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	0.89	1/1184 (0.1%)	1.09	5/1590 (0.3%)
1	B	0.88	1/1183 (0.1%)	1.02	2/1583 (0.1%)
1	C	0.89	1/1199 (0.1%)	1.08	6/1609 (0.4%)
1	D	0.89	1/1192 (0.1%)	1.03	3/1595 (0.2%)
All	All	0.89	4/4758 (0.1%)	1.05	16/6377 (0.3%)

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	A	0	2
1	C	0	2
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	175	LEU	CG-CD1	-6.74	1.30	1.52
1	D	175	LEU	CG-CD1	-6.64	1.30	1.52
1	A	175	LEU	CG-CD2	-6.01	1.32	1.52
1	C	175	LEU	CG-CD2	-5.99	1.32	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	152	ILE	CB-CA-C	7.41	121.67	110.62
1	D	175	LEU	CD1-CG-CD2	-7.28	94.79	110.80
1	C	152	ILE	CB-CA-C	6.97	121.00	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	175	LEU	CD1-CG-CD2	-6.92	95.57	110.80
1	A	111	ASN	N-CA-C	6.67	125.00	110.80
1	A	143	GLY	N-CA-C	6.39	118.47	110.29
1	A	175	LEU	CD1-CG-CD2	-6.28	96.98	110.80
1	C	175	LEU	CD1-CG-CD2	-6.17	97.22	110.80
1	C	163	LEU	CA-CB-CG	5.81	136.63	116.30
1	B	107	LEU	N-CA-C	5.65	117.52	111.36
1	C	110	PRO	CA-C-N	5.63	131.83	121.70
1	C	110	PRO	C-N-CA	5.63	131.83	121.70
1	D	107	LEU	N-CA-C	5.60	117.47	111.36
1	A	163	LEU	CA-CB-CG	5.50	135.55	116.30
1	D	148	VAL	CB-CA-C	5.08	117.60	110.99
1	C	110	PRO	O-C-N	-5.04	116.77	123.01

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	110	PRO	Peptide
1	A	111	ASN	Peptide
1	C	110	PRO	Peptide
1	C	111	ASN	Peptide

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	1162	0	1148	30	0
1	B	1156	0	1151	28	0
1	C	1174	0	1165	31	0
1	D	1162	0	1157	33	0
2	A	3	0	0	0	0
2	B	1	0	0	0	0
2	C	3	0	0	0	0
2	D	1	0	0	0	0
3	A	8	0	12	0	0
3	C	8	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	45	0	0	1	0
4	B	39	0	0	3	0
4	C	42	0	0	2	0
4	D	30	0	0	2	0
All	All	4834	0	4645	121	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 13.

All (121) 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:111:ASN:HA	1:C:112:THR:HG23	1.18	1.15
1:A:111:ASN:HA	1:A:112:THR:CG2	1.80	1.11
1:A:111:ASN:CA	1:A:112:THR:HG23	1.90	1.00
1:A:111:ASN:HA	1:A:112:THR:HG23	0.99	0.98
1:D:114:ASN:C	1:D:116:PRO:HD3	1.91	0.95
1:D:114:ASN:O	1:D:116:PRO:HD3	1.74	0.88
1:C:111:ASN:CA	1:C:112:THR:HG23	2.06	0.84
1:B:120:LYS:HD2	1:B:123:SER:OG	1.76	0.84
1:D:120:LYS:HD2	1:D:123:SER:OG	1.77	0.84
1:A:132:ARG:HH11	1:A:132:ARG:CG	1.93	0.80
1:C:132:ARG:CG	1:C:132:ARG:HH11	1.94	0.80
1:C:110:PRO:CB	1:C:111:ASN:HB2	2.13	0.77
1:B:132:ARG:HH11	1:B:132:ARG:CG	1.98	0.77
1:D:132:ARG:CG	1:D:132:ARG:HH11	1.99	0.75
1:C:111:ASN:HA	1:C:112:THR:CG2	2.09	0.75
1:A:62:GLN:HE21	1:A:67:ARG:HE	1.37	0.72
1:C:62:GLN:HE21	1:C:67:ARG:HE	1.38	0.70
1:A:135:LEU:HD23	1:A:148:VAL:HB	1.76	0.68
1:C:110:PRO:HB3	1:C:111:ASN:HB2	1.75	0.68
1:C:135:LEU:HD23	1:C:148:VAL:HB	1.75	0.68
1:A:132:ARG:HH11	1:A:132:ARG:HG3	1.60	0.66
1:D:72:ALA:HA	1:D:82:GLN:HG3	1.78	0.66
1:D:41:GLU:N	4:D:13:HOH:O	2.30	0.65
1:B:114:ASN:O	1:B:115:LEU:HB2	1.96	0.65
1:B:125:MET:O	4:B:197:HOH:O	2.15	0.64
1:B:62:GLN:HE21	1:B:67:ARG:HE	1.44	0.64
1:C:62:GLN:NE2	1:C:67:ARG:HE	1.97	0.63
1:C:132:ARG:HH11	1:C:132:ARG:HG3	1.64	0.63
1:B:72:ALA:HA	1:B:82:GLN:HG3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:GLN:NE2	1:A:67:ARG:HE	1.97	0.62
1:C:142:ARG:HD2	4:C:207:HOH:O	2.00	0.62
1:D:62:GLN:HE21	1:D:67:ARG:HE	1.45	0.62
1:A:63:ASN:HD22	1:A:63:ASN:C	2.10	0.60
1:B:135:LEU:HD23	1:B:148:VAL:HB	1.83	0.60
1:C:130:TYR:HB3	1:C:153:THR:HG22	1.82	0.60
1:A:130:TYR:HB3	1:A:153:THR:HG22	1.83	0.59
1:C:132:ARG:HH11	1:C:132:ARG:HG2	1.67	0.59
1:B:132:ARG:HH11	1:B:132:ARG:HG2	1.67	0.59
1:B:181:GLU:CD	1:D:120:LYS:HE2	2.27	0.59
1:C:63:ASN:C	1:C:63:ASN:HD22	2.11	0.58
1:D:135:LEU:HD23	1:D:148:VAL:HB	1.86	0.58
1:B:62:GLN:NE2	1:B:67:ARG:HE	2.02	0.57
1:B:62:GLN:HE22	1:B:67:ARG:HH21	1.53	0.57
1:D:62:GLN:NE2	1:D:67:ARG:HE	2.03	0.56
1:D:132:ARG:HH11	1:D:132:ARG:HG2	1.68	0.56
1:B:63:ASN:C	1:B:63:ASN:HD22	2.14	0.56
1:A:70:LYS:HE2	1:A:82:GLN:HG2	1.88	0.56
1:D:62:GLN:HE22	1:D:67:ARG:HH21	1.54	0.55
1:D:63:ASN:HD22	1:D:63:ASN:C	2.15	0.55
1:D:114:ASN:O	1:D:116:PRO:CD	2.53	0.55
1:B:47:LYS:HE3	1:B:177:ASP:OD1	2.07	0.54
1:D:110:PRO:HB2	1:D:111:ASN:HB2	1.88	0.54
1:A:132:ARG:HH11	1:A:132:ARG:HG2	1.70	0.54
1:D:82:GLN:NE2	4:D:33:HOH:O	2.40	0.54
1:D:132:ARG:HH11	1:D:132:ARG:HG3	1.72	0.54
1:C:70:LYS:HE2	1:C:82:GLN:HG2	1.91	0.53
1:B:114:ASN:ND2	4:B:196:HOH:O	2.41	0.53
1:D:139:GLU:HG3	1:D:144:ARG:HG3	1.91	0.53
1:A:152:ILE:O	1:A:155:GLY:HA2	2.10	0.52
1:D:110:PRO:CB	1:D:111:ASN:HB2	2.40	0.52
1:A:62:GLN:HE22	1:A:67:ARG:HH21	1.58	0.51
1:B:132:ARG:HH11	1:B:132:ARG:HG3	1.72	0.51
1:D:47:LYS:HE3	1:D:177:ASP:OD1	2.10	0.51
1:A:47:LYS:HE3	1:A:177:ASP:OD1	2.10	0.51
1:B:139:GLU:HG3	1:B:144:ARG:HG3	1.93	0.51
1:C:47:LYS:HE3	1:C:177:ASP:OD1	2.11	0.51
1:C:62:GLN:HE22	1:C:67:ARG:HH21	1.59	0.50
1:C:152:ILE:O	1:C:155:GLY:HA2	2.11	0.50
1:B:82:GLN:NE2	4:B:33:HOH:O	2.46	0.49
1:A:135:LEU:CD2	1:A:148:VAL:HB	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:ASP:CG	1:A:113:ASP:O	2.56	0.49
1:C:140:ASN:C	1:C:140:ASN:HD22	2.21	0.49
1:D:140:ASN:C	1:D:140:ASN:HD22	2.20	0.48
1:C:135:LEU:CD2	1:C:148:VAL:HB	2.43	0.48
1:B:120:LYS:HD3	1:B:136:ASP:CG	2.39	0.48
1:D:115:LEU:HD11	1:D:140:ASN:HA	1.95	0.48
1:B:140:ASN:HD22	1:B:140:ASN:C	2.22	0.48
1:A:140:ASN:HD22	1:A:140:ASN:C	2.22	0.47
1:A:63:ASN:ND2	1:A:65:ARG:H	2.12	0.47
1:C:63:ASN:ND2	1:C:65:ARG:H	2.13	0.47
1:B:63:ASN:ND2	1:B:65:ARG:H	2.13	0.47
1:D:120:LYS:HD3	1:D:136:ASP:CG	2.39	0.47
1:D:70:LYS:HE2	1:D:82:GLN:HG2	1.97	0.46
1:C:132:ARG:CG	1:C:132:ARG:NH1	2.64	0.46
1:B:70:LYS:HE2	1:B:82:GLN:HG2	1.96	0.46
1:A:72:ALA:HA	1:A:82:GLN:HG3	1.96	0.46
1:C:72:ALA:HA	1:C:82:GLN:HG3	1.98	0.46
1:C:132:ARG:HG3	1:C:132:ARG:NH1	2.31	0.46
1:A:54:LYS:HB3	1:A:74:ILE:O	2.15	0.46
1:C:54:LYS:HB3	1:C:74:ILE:O	2.16	0.45
1:D:63:ASN:ND2	1:D:65:ARG:H	2.13	0.45
1:A:52:GLN:O	1:A:53:SER:OG	2.33	0.45
1:A:111:ASN:N	1:A:112:THR:HG23	2.30	0.44
1:A:110:PRO:CB	1:A:111:ASN:HB2	2.48	0.44
1:C:163:LEU:HD21	1:C:171:PHE:CD2	2.53	0.44
1:B:132:ARG:CG	1:B:132:ARG:NH1	2.67	0.44
1:B:110:PRO:HB2	1:B:111:ASN:HB2	2.00	0.44
1:C:52:GLN:O	1:C:53:SER:OG	2.30	0.44
1:D:62:GLN:NE2	1:D:67:ARG:HH21	2.16	0.44
1:A:132:ARG:HG3	1:A:132:ARG:NH1	2.28	0.43
1:C:110:PRO:CA	1:C:111:ASN:HB2	2.49	0.43
1:C:113:ASP:O	1:C:113:ASP:CG	2.61	0.43
1:B:125:MET:HE2	1:B:134:TYR:HE1	1.84	0.42
1:D:163:LEU:HD21	1:D:171:PHE:CD2	2.54	0.42
1:D:115:LEU:HD13	1:D:115:LEU:O	2.18	0.42
1:A:163:LEU:HD21	1:A:171:PHE:CD2	2.54	0.42
1:D:163:LEU:HD12	1:D:168:MET:HG2	2.02	0.42
1:A:111:ASN:CA	1:A:112:THR:CG2	2.71	0.42
1:C:133:TYR:HA	1:C:149:SER:O	2.20	0.42
1:D:125:MET:HE2	1:D:134:TYR:HE1	1.84	0.42
1:A:133:TYR:HA	1:A:149:SER:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:GLN:NE2	1:B:67:ARG:HH21	2.16	0.41
1:C:52:GLN:NE2	4:C:8:HOH:O	2.53	0.41
1:D:120:LYS:HD2	1:D:123:SER:HG	1.80	0.41
1:B:163:LEU:HD21	1:B:171:PHE:CD2	2.56	0.41
1:D:152:ILE:HD13	1:D:157:PRO:HG2	2.03	0.41
1:A:62:GLN:NE2	1:A:67:ARG:HH21	2.18	0.41
1:B:152:ILE:HD13	1:B:157:PRO:HG2	2.03	0.41
1:A:41:GLU:N	4:A:203:HOH:O	2.54	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	144/146 (99%)	135 (94%)	6 (4%)	3 (2%)	5	2
1	B	138/146 (94%)	133 (96%)	3 (2%)	2 (1%)	9	5
1	C	145/146 (99%)	135 (93%)	7 (5%)	3 (2%)	5	2
1	D	139/146 (95%)	135 (97%)	3 (2%)	1 (1%)	18	15
All	All	566/584 (97%)	538 (95%)	19 (3%)	9 (2%)	7	4

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	113	ASP
1	B	115	LEU
1	C	111	ASN
1	C	112	THR
1	C	113	ASP
1	A	112	THR
1	B	116	PRO

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Mol	Chain	Res	Type
1	D	115	LEU
1	A	111	ASN

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	121/124 (98%)	109 (90%)	12 (10%)	7	5
1	B	121/124 (98%)	112 (93%)	9 (7%)	13	10
1	C	123/124 (99%)	108 (88%)	15 (12%)	5	2
1	D	122/124 (98%)	113 (93%)	9 (7%)	13	10
All	All	487/496 (98%)	442 (91%)	45 (9%)	8	6

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	LEU
1	A	54	LYS
1	A	63	ASN
1	A	112	THR
1	A	113	ASP
1	A	121	LEU
1	A	132	ARG
1	A	140	ASN
1	A	148	VAL
1	A	152	ILE
1	A	177	ASP
1	A	185	ASN
1	B	49	LEU
1	B	63	ASN
1	B	115	LEU
1	B	132	ARG
1	B	140	ASN
1	B	142	ARG
1	B	148	VAL

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Mol	Chain	Res	Type
1	B	175	LEU
1	B	177	ASP
1	C	49	LEU
1	C	54	LYS
1	C	63	ASN
1	C	111	ASN
1	C	112	THR
1	C	113	ASP
1	C	121	LEU
1	C	132	ARG
1	C	140	ASN
1	C	148	VAL
1	C	152	ILE
1	C	177	ASP
1	C	180[A]	GLU
1	C	180[B]	GLU
1	C	185	ASN
1	D	49	LEU
1	D	63	ASN
1	D	115	LEU
1	D	132	ARG
1	D	140	ASN
1	D	142	ARG
1	D	148	VAL
1	D	175	LEU
1	D	177	ASP

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

Mol	Chain	Res	Type
1	A	42	GLN
1	A	62	GLN
1	A	63	ASN
1	A	160	GLN
1	B	63	ASN
1	B	111	ASN
1	B	140	ASN
1	C	42	GLN
1	C	62	GLN
1	C	63	ASN
1	C	140	ASN
1	D	42	GLN

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Mol	Chain	Res	Type
1	D	62	GLN
1	D	63	ASN
1	D	140	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	$\# Z > 2$
3	EDO	A	3318	-	3,3,3	0.44	0	2,2,2	0.50	0
3	EDO	C	188	-	3,3,3	0.66	0	2,2,2	0.28	0
3	EDO	A	187	-	3,3,3	0.65	0	2,2,2	0.16	0
3	EDO	C	3318	-	3,3,3	0.52	0	2,2,2	0.40	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
3	EDO	A	3318	-	-	0/1/1/1	-
3	EDO	C	188	-	-	1/1/1/1	-
3	EDO	A	187	-	-	1/1/1/1	-
3	EDO	C	3318	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	188	EDO	O1-C1-C2-O2
3	A	187	EDO	O1-C1-C2-O2

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	145/146 (99%)	-1.32	0 100 100	15, 33, 57, 67	1 (0%)
1	B	140/146 (95%)	-1.27	0 100 100	15, 33, 55, 88	4 (2%)
1	C	145/146 (99%)	-1.32	0 100 100	15, 33, 57, 67	2 (1%)
1	D	140/146 (95%)	-1.32	0 100 100	15, 33, 55, 99	5 (3%)
All	All	570/584 (97%)	-1.31	0 100 100	15, 33, 57, 99	12 (2%)

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
3	EDO	A	3318	4/4	0.98	0.07	28,28,29,31	4
3	EDO	A	187	4/4	0.98	0.04	39,40,40,41	0
3	EDO	C	3318	4/4	0.98	0.07	26,26,27,27	4
3	EDO	C	188	4/4	0.98	0.04	41,42,43,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CL	C	187	1/1	0.99	0.09	49,49,49,49	0
2	CL	A	1	1/1	0.99	0.02	30,30,30,30	0
2	CL	A	186	1/1	0.99	0.05	48,48,48,48	0
2	CL	A	188	1/1	0.99	0.03	45,45,45,45	0
2	CL	B	1	1/1	0.99	0.02	48,48,48,48	0
2	CL	C	1	1/1	1.00	0.02	47,47,47,47	0
2	CL	D	1	1/1	1.00	0.02	48,48,48,48	0
2	CL	C	186	1/1	1.00	0.02	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.