



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:35 PM GMT

PDB ID : 4CNT
Title : CRYSTAL STRUCTURE OF WT HUMAN CRMP-4
Authors : Ponnusamy, R.; Lohkamp, B.
Deposited on : 2014-01-24
Resolution : 2.65 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

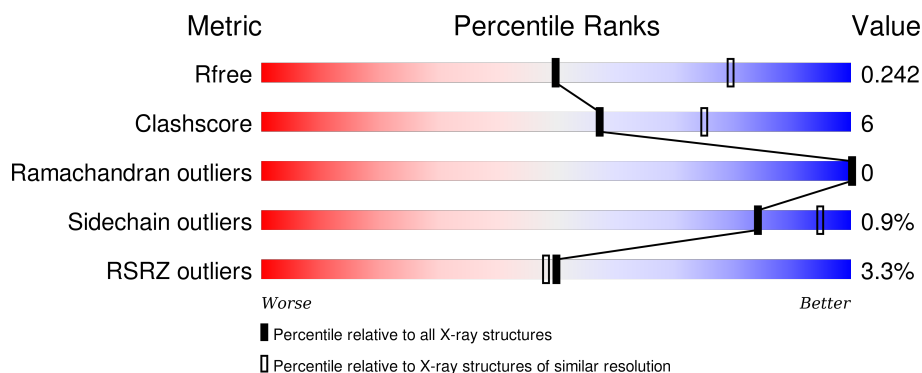
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	91344	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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. 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	570	<div> <div> <div></div> <div>74%</div> <div>11%</div> <div>•</div> <div>15%</div> </div> </div>
1	B	570	<div> <div> <div>2%</div> <div></div> <div>74%</div> <div>11%</div> <div>•</div> <div>15%</div> </div> </div>
1	C	570	<div> <div> <div>5%</div> <div></div> <div>74%</div> <div>10%</div> <div>•</div> <div>15%</div> </div> </div>
1	D	570	<div> <div> <div>4%</div> <div></div> <div>74%</div> <div>11%</div> <div></div> <div>15%</div> </div> </div>

2 Entry composition [i](#)

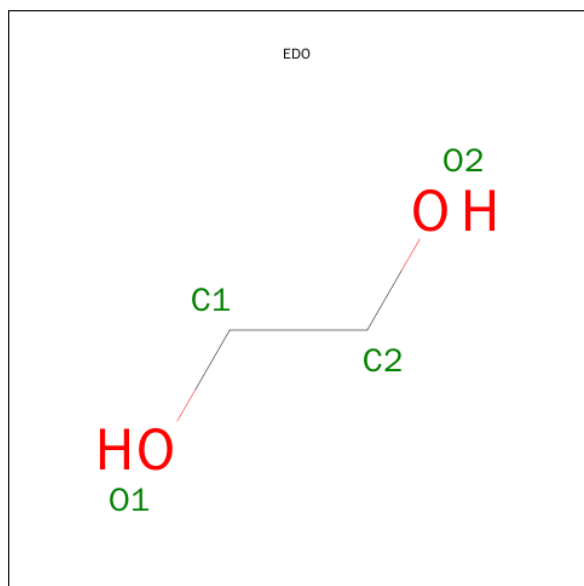
There are 4 unique types of molecules in this entry. The entry contains 14948 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DIHYDROPYRIMIDINASE-LIKE 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	485	Total	C	N	O	S	0	0	0
			3732	2357	632	721	22			
1	B	486	Total	C	N	O	S	0	0	0
			3736	2359	633	722	22			
1	C	485	Total	C	N	O	S	0	0	0
			3732	2357	632	721	22			
1	D	485	Total	C	N	O	S	0	0	0
			3732	2357	632	721	22			

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

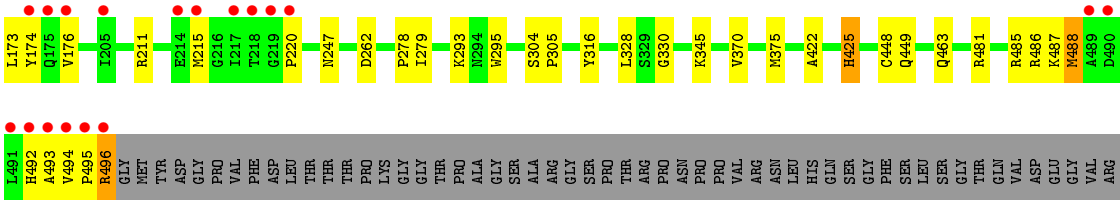
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total Na 1 1	0	0

- Molecule 4 is water.

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

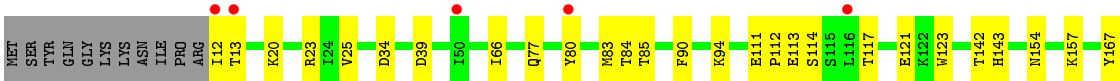
• Molecule 1: DIHYDROPYRIMIDINASE-LIKE 3





SER
ALA
SER
LYS
ARG
ILE
VAL
ALA
PRO
PRO
GLY
ASP
PRO
GLY
ARG
VAL
PHE
SER
ASN
ILE
THR
SER
LEU
SER

• Molecule 1: DIHYDROPYRIMIDINASE-LIKE 3



GLN
SER
GLY
PHE
SER
LEU
SER
GLY
THR
GLN
VAL
ASP
GLU
GLY
VAL
ARG
SER
ALA
SER
LYS
ARG
ILE
VAL
ALA
PRO
PRO
GLY
GLY
ARG
PHE
SER
ASN
ILE
THR
SER
LEU
SER

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.37Å 157.65Å 86.38Å 90.00° 113.03° 90.00°	Depositor
Resolution (Å)	55.97 – 2.65 55.97 – 2.66	Depositor EDS
% Data completeness (in resolution range)	99.8 (55.97-2.65) 99.9 (55.97-2.66)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.187 , 0.236 0.194 , 0.242	Depositor DCC
R_{free} test set	3105 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.465	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.6	EDS
Estimated twinning fraction	0.685 for H, K, L 0.315 for L, -K, H 0.328 for l,-k,h	Xtriage
Reported twinning fraction	0.685 for H, K, L 0.315 for L, -K, H	Depositor
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	1 of 61244 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	14948	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: NA, 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	0.84	1/3810 (0.0%)	0.83	0/5168
1	B	0.76	0/3814	0.84	2/5173 (0.0%)
1	C	0.76	0/3810	0.83	2/5168 (0.0%)
1	D	0.72	0/3810	0.81	1/5168 (0.0%)
All	All	0.77	1/15244 (0.0%)	0.83	5/20677 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	76	TYR	CE1-CZ	-5.28	1.31	1.38

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	34	ASP	CB-CG-OD2	7.02	124.62	118.30
1	C	39	ASP	CB-CG-OD2	-5.71	113.16	118.30
1	C	23	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	B	165	MET	CG-SD-CE	5.41	108.85	100.20
1	D	415	ASP	CB-CG-OD1	5.34	123.10	118.30

There are no chirality outliers.

There are no planarity outliers.

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	3732	0	3679	43	0
1	B	3736	0	3682	48	0
1	C	3732	0	3679	41	0
1	D	3732	0	3679	51	0
2	A	4	0	6	0	0
3	D	1	0	0	0	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
4	C	3	0	0	0	0
4	D	2	0	0	0	0
All	All	14948	0	14725	178	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 (178) 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:B:111:GLU:O	1:B:114:SER:HB3	1.81	0.80
1:D:430:GLU:N	1:D:430:GLU:OE1	2.14	0.79
1:C:493:ALA:O	1:C:495:PRO:HD3	1.81	0.79
1:A:12:ILE:HG22	1:A:12:ILE:O	1.85	0.77
1:B:12:ILE:O	1:B:12:ILE:HG22	1.84	0.76
1:C:80:TYR:O	1:C:83:MET:HB2	1.88	0.74
1:C:12:ILE:HG22	1:C:12:ILE:O	1.86	0.74
1:D:422:ALA:HA	1:D:425:HIS:ND1	2.01	0.74
1:C:481:ARG:O	1:C:485:ARG:HG3	1.89	0.73
1:D:113:GLU:H	1:D:143:HIS:CE1	2.04	0.73
1:A:486:ARG:O	1:A:489:ALA:HB3	1.89	0.73
1:A:481:ARG:O	1:A:485:ARG:HG3	1.93	0.69
1:D:476:ASP:HB3	1:D:480:LYS:CE	2.25	0.66
1:C:166:VAL:HG12	1:C:167:TYR:N	2.11	0.65
1:D:262:ASP:OD1	1:D:316:TYR:OH	2.15	0.65
1:D:23:ARG:HG3	1:D:23:ARG:HH11	1.63	0.64
1:D:142:THR:HB	1:D:174:TYR:CD1	2.33	0.64
1:D:476:ASP:HB3	1:D:480:LYS:HE3	1.80	0.63
1:A:491:LEU:HD12	1:B:215:MET:HG3	1.79	0.63
1:B:80:TYR:CE1	1:B:81:LYS:HG3	2.33	0.63
1:C:293:LYS:O	1:D:496:ARG:HG2	1.98	0.63
1:C:262:ASP:OD1	1:C:316:TYR:OH	2.17	0.62
1:C:142:THR:HB	1:C:174:TYR:CD1	2.36	0.61
1:A:113:GLU:H	1:A:143:HIS:CE1	2.20	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:ARG:HA	1:A:37:MET:O	2.02	0.60
1:D:111:GLU:O	1:D:114:SER:HB3	2.01	0.60
1:B:422:ALA:HA	1:B:425:HIS:CE1	2.37	0.60
1:D:84:THR:HG22	1:D:85:THR:O	2.02	0.60
1:B:422:ALA:HA	1:B:425:HIS:ND1	2.16	0.60
1:B:90:PHE:CE1	1:B:94:LYS:HE2	2.36	0.60
1:B:12:ILE:O	1:B:12:ILE:CG2	2.50	0.60
1:D:170:TYR:HB3	1:D:173:LEU:HB3	1.84	0.60
1:B:262:ASP:OD1	1:B:316:TYR:OH	2.19	0.60
1:A:84:THR:HG22	1:A:85:THR:O	2.01	0.59
1:A:12:ILE:CG2	1:A:12:ILE:O	2.51	0.59
1:A:111:GLU:O	1:A:114:SER:HB3	2.03	0.59
1:C:113:GLU:H	1:C:143:HIS:CE1	2.20	0.59
1:A:51:VAL:HG12	1:A:52:PRO:O	2.03	0.58
1:B:334:CYS:HB3	1:B:354:GLY:HA3	1.85	0.58
1:C:12:ILE:O	1:C:12:ILE:CG2	2.51	0.58
1:D:333:HIS:CE1	1:D:356:ASN:HA	2.39	0.57
1:B:23:ARG:NH1	1:B:377:GLU:OE2	2.33	0.57
1:C:494:VAL:HG11	1:C:496:ARG:NH1	2.19	0.57
1:A:262:ASP:OD1	1:A:316:TYR:OH	2.18	0.56
1:A:113:GLU:HA	1:A:113:GLU:OE1	2.05	0.56
1:D:484:ALA:O	1:D:488:MET:HB2	2.05	0.56
1:B:295:TRP:CZ2	1:B:345:LYS:HA	2.41	0.56
1:B:80:TYR:CE1	1:B:81:LYS:CG	2.89	0.56
1:D:485:ARG:HA	1:D:488:MET:HE2	1.87	0.55
1:B:476:ASP:CG	1:B:480:LYS:HE3	2.27	0.55
1:B:84:THR:HG22	1:B:85:THR:O	2.07	0.55
1:D:295:TRP:CZ2	1:D:345:LYS:HA	2.42	0.55
1:D:422:ALA:HA	1:D:425:HIS:CE1	2.41	0.55
1:D:113:GLU:N	1:D:143:HIS:CE1	2.74	0.54
1:A:12:ILE:HG22	1:A:39:ASP:O	2.08	0.54
1:D:171:LYS:O	1:D:172:ASP:HB2	2.08	0.54
1:D:142:THR:HB	1:D:174:TYR:HD1	1.72	0.53
1:A:485:ARG:O	1:A:489:ALA:HB2	2.08	0.53
1:D:421:SER:HB2	1:D:435:GLU:OE2	2.07	0.53
1:D:476:ASP:HB3	1:D:480:LYS:HE2	1.91	0.53
1:B:171:LYS:O	1:B:172:ASP:HB2	2.07	0.52
1:C:220:PRO:HD2	1:D:496:ARG:NH1	2.24	0.52
1:B:167:TYR:O	1:B:176:VAL:HG22	2.10	0.52
1:D:154:ASN:HA	1:D:157:LYS:HG2	1.93	0.51
1:C:171:LYS:O	1:C:172:ASP:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:CYS:HB3	1:A:354:GLY:HA3	1.93	0.51
1:D:154:ASN:OD1	1:D:157:LYS:HE2	2.11	0.51
1:A:167:TYR:O	1:A:176:VAL:HG22	2.11	0.50
1:C:80:TYR:O	1:C:83:MET:N	2.40	0.50
1:D:23:ARG:HG3	1:D:23:ARG:NH1	2.25	0.50
1:D:80:TYR:O	1:D:83:MET:HB2	2.11	0.50
1:B:80:TYR:HE1	1:B:81:LYS:HG3	1.76	0.50
1:D:112:PRO:HA	1:D:143:HIS:ND1	2.27	0.50
1:C:167:TYR:O	1:C:176:VAL:HG22	2.12	0.49
1:D:167:TYR:O	1:D:176:VAL:HG22	2.11	0.49
1:D:90:PHE:CE2	1:D:94:LYS:HE2	2.46	0.49
1:C:304:SER:HA	1:C:305:PRO:C	2.31	0.49
1:B:484:ALA:O	1:B:488:MET:HB2	2.13	0.49
1:C:114:SER:OG	1:C:115:SER:N	2.45	0.49
1:C:112:PRO:O	1:C:113:GLU:HB2	2.12	0.49
1:D:421:SER:HA	1:D:435:GLU:OE2	2.13	0.49
1:A:171:LYS:O	1:A:172:ASP:HB2	2.13	0.48
1:C:422:ALA:HA	1:C:425:HIS:CE1	2.49	0.48
1:C:84:THR:HG22	1:C:85:THR:O	2.13	0.48
1:A:295:TRP:CZ2	1:A:345:LYS:HA	2.49	0.48
1:C:166:VAL:HG12	1:C:167:TYR:H	1.78	0.48
1:A:16:ARG:CB	1:A:36:TYR:OH	2.62	0.48
1:D:333:HIS:ND1	1:D:356:ASN:HA	2.29	0.48
1:A:484:ALA:O	1:A:488:MET:HG3	2.14	0.48
1:C:173:LEU:HG	1:C:174:TYR:N	2.28	0.47
1:A:52:PRO:O	1:A:55:VAL:HG23	2.15	0.47
1:C:494:VAL:CG1	1:C:496:ARG:NH1	2.78	0.47
1:A:16:ARG:HB3	1:A:36:TYR:OH	2.14	0.47
1:A:51:VAL:CG1	1:A:55:VAL:HG21	2.46	0.46
1:D:77:GLN:HA	1:D:84:THR:HG23	1.97	0.46
1:A:333:HIS:CE1	1:A:358:VAL:HG12	2.50	0.46
1:A:419:ILE:HG13	1:A:438:GLU:HG3	1.97	0.46
1:D:448:CYS:O	1:D:449:GLN:HB2	2.16	0.46
1:B:476:ASP:CB	1:B:480:LYS:HE3	2.46	0.46
1:D:300:ALA:O	1:D:355:THR:HG22	2.16	0.46
1:B:476:ASP:HB3	1:B:480:LYS:HE3	1.98	0.46
1:B:485:ARG:HA	1:B:488:MET:HE2	1.98	0.46
1:B:165:MET:CE	1:B:198:HIS:HB2	2.46	0.46
1:C:166:VAL:CG1	1:C:167:TYR:N	2.77	0.46
1:C:448:CYS:O	1:C:449:GLN:HB2	2.16	0.46
1:B:90:PHE:CE2	1:B:463:GLN:HB2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:496:ARG:NH2	1:D:220:PRO:HD2	2.31	0.45
1:C:23:ARG:HB2	1:C:32:TYR:CE1	2.51	0.45
1:A:76:TYR:CE1	1:A:92:GLY:HA3	2.51	0.45
1:B:80:TYR:CD1	1:B:80:TYR:C	2.86	0.45
1:D:304:SER:HA	1:D:305:PRO:C	2.37	0.45
1:B:142:THR:HB	1:B:174:TYR:CD1	2.52	0.45
1:B:304:SER:HA	1:B:305:PRO:C	2.37	0.45
1:D:77:GLN:HA	1:D:84:THR:CG2	2.47	0.44
1:B:270:LYS:HB3	1:B:270:LYS:HE3	1.66	0.44
1:A:492:HIS:ND1	1:A:492:HIS:N	2.62	0.44
1:B:113:GLU:H	1:B:143:HIS:CE1	2.35	0.44
1:C:278:PRO:HD2	1:C:328:LEU:O	2.18	0.44
1:A:448:CYS:O	1:A:449:GLN:HB2	2.17	0.44
1:C:295:TRP:CZ2	1:C:345:LYS:HA	2.52	0.44
1:D:173:LEU:HD21	1:D:174:TYR:CE2	2.53	0.44
1:B:448:CYS:O	1:B:449:GLN:HB2	2.18	0.44
1:B:286:ASP:OD2	1:B:311:PRO:HA	2.18	0.44
1:A:77:GLN:HA	1:A:84:THR:CG2	2.48	0.43
1:A:51:VAL:HG13	1:A:52:PRO:HD2	1.99	0.43
1:D:117:THR:O	1:D:121:GLU:HG2	2.18	0.43
1:A:279:ILE:HA	1:A:330:GLY:O	2.18	0.43
1:B:112:PRO:O	1:B:113:GLU:HB2	2.17	0.43
1:A:300:ALA:O	1:A:355:THR:HG22	2.19	0.43
1:D:25:VAL:HB	1:D:66:ILE:HG22	1.99	0.43
1:B:300:ALA:O	1:B:355:THR:HG22	2.18	0.43
1:A:278:PRO:HD2	1:A:328:LEU:O	2.19	0.43
1:B:20:LYS:HA	1:B:34:ASP:OD1	2.19	0.43
1:B:165:MET:HE2	1:B:198:HIS:HB2	2.00	0.43
1:A:77:GLN:HA	1:A:84:THR:HG23	2.01	0.43
1:B:16:ARG:HG2	1:B:36:TYR:OH	2.19	0.43
1:D:425:HIS:C	1:D:425:HIS:CD2	2.92	0.43
1:B:117:THR:O	1:B:121:GLU:HG2	2.19	0.43
1:D:20:LYS:HA	1:D:34:ASP:OD1	2.19	0.42
1:A:142:THR:HB	1:A:174:TYR:CD1	2.53	0.42
1:D:278:PRO:HD2	1:D:328:LEU:O	2.19	0.42
1:C:20:LYS:HA	1:C:34:ASP:OD1	2.18	0.42
1:A:304:SER:HA	1:A:305:PRO:C	2.39	0.42
1:D:12:ILE:O	1:D:13:THR:C	2.58	0.42
1:C:496:ARG:NH2	1:D:221:GLU:OE1	2.50	0.42
1:C:279:ILE:HA	1:C:330:GLY:O	2.20	0.42
1:C:117:THR:O	1:C:121:GLU:HG2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:25:VAL:HB	1:C:66:ILE:HG22	2.00	0.42
1:D:419:ILE:HG13	1:D:438:GLU:HG3	2.01	0.42
1:A:117:THR:O	1:A:121:GLU:HG2	2.19	0.42
1:B:425:HIS:CD2	1:B:427:SER:HB3	2.54	0.42
1:A:155:LEU:HD23	1:A:159:LYS:HG3	2.01	0.42
1:B:279:ILE:HA	1:B:330:GLY:O	2.20	0.42
1:A:113:GLU:CA	1:A:113:GLU:OE1	2.68	0.42
1:C:247:ASN:ND2	1:C:486:ARG:HD3	2.34	0.42
1:A:20:LYS:HA	1:A:34:ASP:OD1	2.19	0.42
1:B:277:GLU:O	1:B:277:GLU:HG2	2.20	0.42
1:A:25:VAL:HB	1:A:66:ILE:HG22	2.01	0.42
1:D:279:ILE:HA	1:D:330:GLY:O	2.20	0.42
1:C:43:LYS:HD3	1:C:43:LYS:HA	1.77	0.41
1:B:278:PRO:HD2	1:B:328:LEU:O	2.20	0.41
1:B:80:TYR:CZ	1:B:81:LYS:HD2	2.54	0.41
1:B:80:TYR:CD1	1:B:81:LYS:N	2.89	0.41
1:B:130:LYS:HA	1:B:463:GLN:HG3	2.02	0.41
1:A:422:ALA:HA	1:A:425:HIS:CE1	2.55	0.41
1:B:80:TYR:CE1	1:B:81:LYS:HB2	2.55	0.41
1:C:485:ARG:HA	1:C:488:MET:HE2	2.02	0.41
1:D:112:PRO:C	1:D:113:GLU:HG2	2.41	0.41
1:C:130:LYS:HA	1:C:463:GLN:HG3	2.03	0.41
1:B:481:ARG:O	1:B:485:ARG:HG3	2.21	0.40
1:C:492:HIS:H	1:C:492:HIS:CD2	2.38	0.40
1:A:90:PHE:CE2	1:A:94:LYS:HE2	2.55	0.40
1:B:25:VAL:HB	1:B:66:ILE:HG22	2.04	0.40
1:D:12:ILE:HG23	1:D:39:ASP:O	2.20	0.40
1:D:80:TYR:O	1:D:83:MET:N	2.43	0.40
1:B:78:MET:HB3	1:B:78:MET:HE2	1.96	0.40
1:C:211:ARG:HG2	1:C:215:MET:HE2	2.02	0.40
1:C:370:VAL:HG22	1:C:375:MET:HG3	2.04	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	483/570 (85%)	472 (98%)	11 (2%)	0	100	100
1	B	484/570 (85%)	475 (98%)	9 (2%)	0	100	100
1	C	483/570 (85%)	474 (98%)	9 (2%)	0	100	100
1	D	483/570 (85%)	474 (98%)	9 (2%)	0	100	100
All	All	1933/2280 (85%)	1895 (98%)	38 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	402/471 (85%)	398 (99%)	4 (1%)	82	94
1	B	402/471 (85%)	399 (99%)	3 (1%)	88	96
1	C	402/471 (85%)	397 (99%)	5 (1%)	78	92
1	D	402/471 (85%)	400 (100%)	2 (0%)	92	98
All	All	1608/1884 (85%)	1594 (99%)	14 (1%)	84	95

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ARG
1	A	123	TRP
1	A	425	HIS
1	A	492	HIS
1	B	123	TRP
1	B	335	THR
1	B	425	HIS
1	C	123	TRP
1	C	425	HIS
1	C	487	LYS

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Mol	Chain	Res	Type
1	C	488	MET
1	C	496	ARG
1	D	123	TRP
1	D	425	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	424	ASN
1	B	207	GLN
1	C	247	ASN
1	C	492	HIS

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, 1 is monoatomic - leaving 1 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	$\# Z > 2$
2	EDO	A	1497	-	3,3,3	0.92	0	2,2,2	0.06	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	EDO	A	1497	-	-	0/1/1/1	0/0/0/0

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 ⓘ

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	485/570 (85%)	0.09	3 (0%) 90 91	21, 39, 70, 120	0
1	B	486/570 (85%)	0.26	11 (2%) 64 62	22, 49, 76, 94	0
1	C	485/570 (85%)	0.30	26 (5%) 29 27	22, 44, 80, 123	0
1	D	485/570 (85%)	0.37	25 (5%) 31 28	29, 49, 79, 101	0
All	All	1941/2280 (85%)	0.25	65 (3%) 50 48	21, 46, 77, 123	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	495	PRO	6.7
1	C	493	ALA	6.7
1	C	494	VAL	6.6
1	D	493	ALA	6.5
1	D	496	ARG	6.2
1	D	491	LEU	5.6
1	C	12	ILE	5.4
1	C	80	TYR	5.3
1	C	496	ARG	5.0
1	D	492	HIS	5.0
1	C	174	TYR	4.5
1	D	489	ALA	4.1
1	B	12	ILE	4.1
1	D	218	THR	3.9
1	C	170	TYR	3.7
1	C	219	GLY	3.5
1	D	495	PRO	3.5
1	B	293	LYS	3.3
1	D	220	PRO	3.2
1	D	169	ALA	3.2
1	C	175	GLN	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	220	PRO	2.9
1	C	176	VAL	2.9
1	D	174	TYR	2.9
1	C	218	THR	2.9
1	C	172	ASP	2.8
1	D	490	ASP	2.8
1	D	80	TYR	2.8
1	D	291	TRP	2.8
1	B	436	GLY	2.7
1	D	292	SER	2.7
1	C	214	GLU	2.7
1	C	81	LYS	2.7
1	B	444	LEU	2.7
1	C	492	HIS	2.6
1	D	50	ILE	2.6
1	C	217	ILE	2.5
1	C	205	ILE	2.5
1	D	217	ILE	2.5
1	C	171	LYS	2.5
1	C	490	ASP	2.4
1	B	458	ASN	2.4
1	D	494	VAL	2.3
1	D	347	ASN	2.3
1	A	15	ASP	2.3
1	B	420	VAL	2.3
1	B	497	GLY	2.3
1	B	90	PHE	2.3
1	A	490	ASP	2.3
1	D	168	MET	2.2
1	B	220	PRO	2.2
1	C	215	MET	2.2
1	D	13	THR	2.2
1	D	338	THR	2.2
1	D	116	LEU	2.1
1	C	169	ALA	2.1
1	D	12	ILE	2.1
1	D	175	GLN	2.1
1	C	489	ALA	2.1
1	C	13	THR	2.1
1	C	491	LEU	2.1
1	B	456	ASP	2.1
1	D	172	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	13	THR	2.0
1	B	129	GLY	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	NA	D	1497	1/1	0.85	0.13	-1.53	36,36,36,36	0
2	EDO	A	1497	4/4	0.86	0.17	-	27,31,31,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.