



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:34 AM GMT

PDB ID : 3ITK
Title : Crystal structure of human UDP-glucose dehydrogenase Thr131Ala, apo form.
Authors : Chaikuad, A.; Egger, S.; Yue, W.W.; Sethi, R.; Filippakopoulos, P.; Muniz, J.R.C.; von Delft, F.; Bountra, C.; Arrowsmith, C.H.; Weigelt, J.; Edwards, A.M.; Kavanagh, K.L.; Nidetzky, B.; Oppermann, U.; Structural Genomics Consortium (SGC)
Deposited on : 2009-08-28
Resolution : 2.40 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

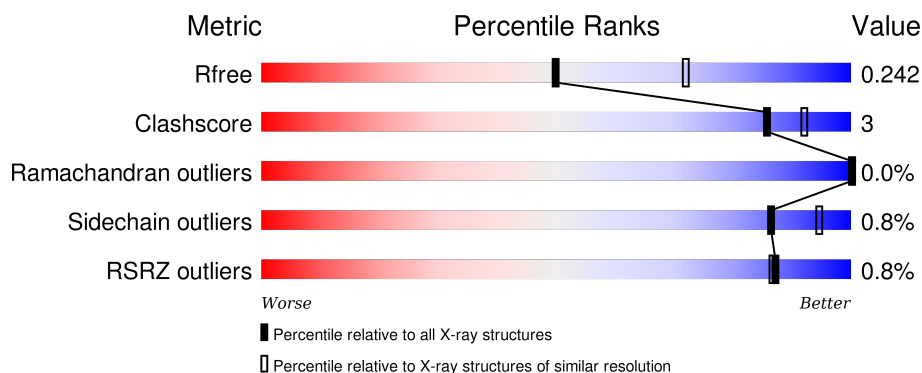
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	467	<div> <div>94%</div> <div>5%</div> </div>
1	B	467	<div> <div>93%</div> <div>5%</div> </div>
1	C	467	<div> <div>94%</div> <div>6%</div> </div>
1	D	467	<div> <div>93%</div> <div>6%</div> </div>
1	E	467	<div> <div>93%</div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	467	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PG4	A	467	-	-	-	X
2	PG4	C	467	-	-	-	X
2	PG4	D	467	-	-	-	X
2	PG4	F	467	-	-	-	X
3	EDO	B	469	-	-	-	X
3	EDO	D	468	-	-	-	X
3	EDO	D	470	-	-	-	X
3	EDO	F	468	-	-	-	X
3	EDO	F	469	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 22977 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.

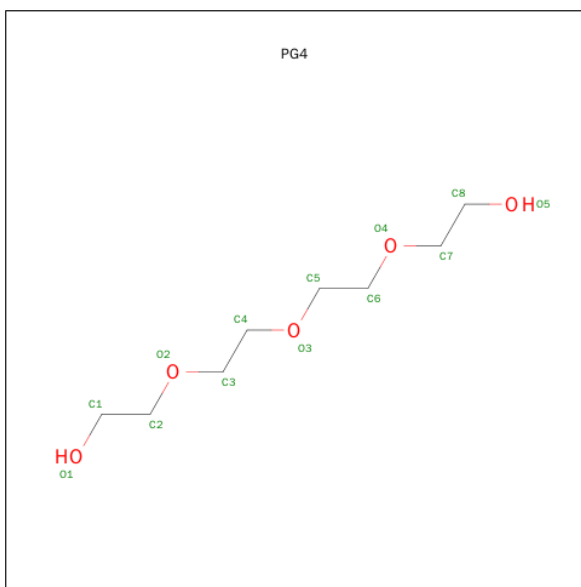
- Molecule 1 is a protein called UDP-glucose 6-dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3591	2277	615	680	19			
1	B	462	Total	C	N	O	S	0	2	0
			3586	2270	615	681	20			
1	C	467	Total	C	N	O	S	0	1	0
			3629	2297	625	688	19			
1	D	467	Total	C	N	O	S	0	0	0
			3622	2295	622	685	20			
1	E	466	Total	C	N	O	S	0	2	0
			3604	2284	619	682	19			
1	F	467	Total	C	N	O	S	0	0	0
			3624	2290	625	690	19			

There are 12 discrepancies between the modelled and reference sequences:

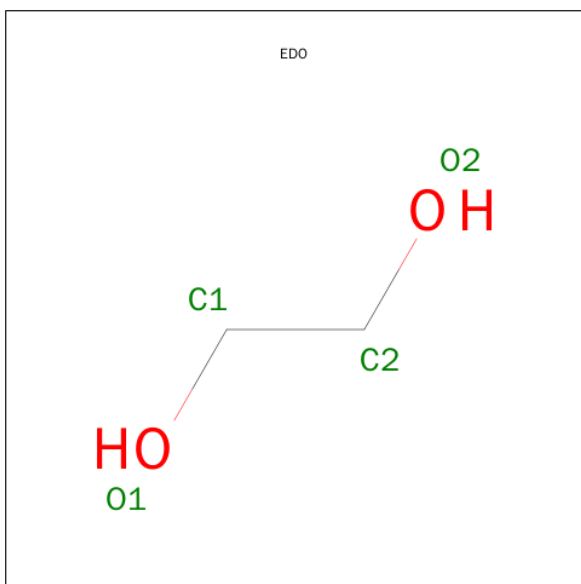
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	EXPRESSION TAG	UNP O60701
A	131	ALA	THR	ENGINEERED	UNP O60701
B	0	SER	-	EXPRESSION TAG	UNP O60701
B	131	ALA	THR	ENGINEERED	UNP O60701
C	0	SER	-	EXPRESSION TAG	UNP O60701
C	131	ALA	THR	ENGINEERED	UNP O60701
D	0	SER	-	EXPRESSION TAG	UNP O60701
D	131	ALA	THR	ENGINEERED	UNP O60701
E	0	SER	-	EXPRESSION TAG	UNP O60701
E	131	ALA	THR	ENGINEERED	UNP O60701
F	0	SER	-	EXPRESSION TAG	UNP O60701
F	131	ALA	THR	ENGINEERED	UNP O60701

- Molecule 2 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	8	5		
2	C	1	Total	C	O	0	0
			13	8	5		
2	C	1	Total	C	O	0	0
			13	8	5		
2	D	1	Total	C	O	0	0
			13	8	5		
2	F	1	Total	C	O	0	0
			13	8	5		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0
3	E	1	Total C O 4 2 2	0	0
3	E	1	Total C O 4 2 2	0	0
3	E	1	Total C O 4 2 2	0	0
3	F	1	Total C O 4 2 2	0	0
3	F	1	Total C O 4 2 2	0	0
3	F	1	Total C O 4 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	211	Total O 211 211	0	0
4	B	171	Total O 171 171	0	0
4	C	215	Total O 215 215	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	202	Total 202	O 202	0	0
4	E	174	Total 174	O 174	0	0
4	F	215	Total 215	O 215	0	0

3 Residue-property plots [i](#)

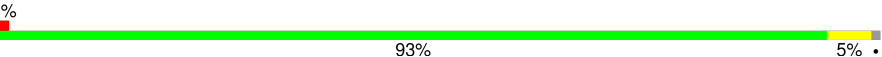
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: UDP-glucose 6-dehydrogenase

Chain A: 

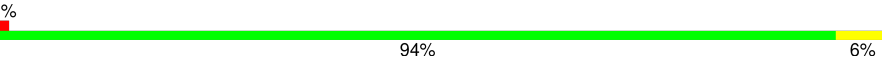


- Molecule 1: UDP-glucose 6-dehydrogenase

Chain B: 



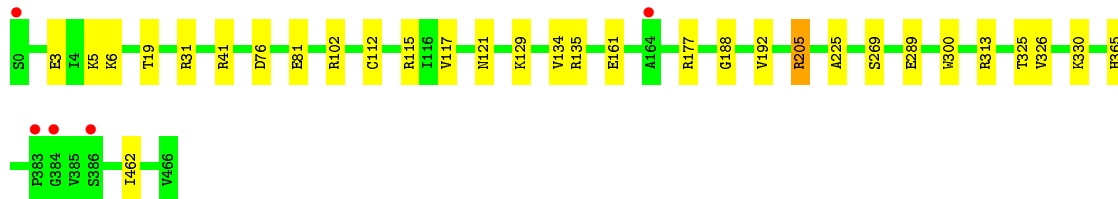
- Molecule 1: UDP-glucose 6-dehydrogenase

Chain C: 

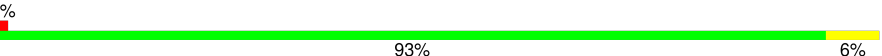


- Molecule 1: UDP-glucose 6-dehydrogenase

Chain D: 



- Molecule 1: UDP-glucose 6-dehydrogenase

Chain E: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.06Å 106.61Å 349.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	52.90 – 2.40 50.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (52.90-2.40) 99.7 (50.98-2.40)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0089	Depositor
R, R_{free}	0.198 , 0.241 0.199 , 0.242	Depositor DCC
R_{free} test set	2134 reflections (1.66%)	DCC
Wilson B-factor (Å ²)	32.0	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 130411 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	22977	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.80% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: PG4, 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	0.73	1/3657 (0.0%)	0.77	8/4956 (0.2%)
1	B	0.73	2/3656 (0.1%)	1.08	8/4953 (0.2%)
1	C	0.75	3/3699 (0.1%)	0.75	7/5012 (0.1%)
1	D	0.75	5/3689 (0.1%)	0.84	7/4998 (0.1%)
1	E	0.71	2/3676 (0.1%)	0.76	9/4984 (0.2%)
1	F	0.76	0/3691	0.79	3/5003 (0.1%)
All	All	0.74	13/22068 (0.1%)	0.84	42/29906 (0.1%)

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	B	0	1
1	D	0	1
1	E	0	1
All	All	0	3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	31	ARG	CZ-NH1	-12.51	1.16	1.33
1	D	121	ASN	CG-ND2	-7.02	1.15	1.32
1	C	31	ARG	CZ-NH1	-6.81	1.24	1.33
1	A	31	ARG	CZ-NH1	-6.57	1.24	1.33
1	E	117	VAL	CB-CG2	-6.51	1.39	1.52
1	E	31	ARG	CZ-NH1	-6.31	1.24	1.33
1	D	31	ARG	CZ-NH1	-6.30	1.24	1.33
1	D	121	ASN	CG-OD1	-6.25	1.10	1.24
1	C	112	CYS	CB-SG	-6.09	1.71	1.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	112	CYS	CB-SG	-5.91	1.72	1.81
1	B	117	VAL	CB-CG2	-5.49	1.41	1.52
1	D	117	VAL	CB-CG2	-5.33	1.41	1.52
1	C	187	GLU	CB-CG	5.17	1.61	1.52

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	31	ARG	NE-CZ-NH2	39.92	140.26	120.30
1	B	31	ARG	NE-CZ-NH1	-33.93	103.33	120.30
1	D	205	ARG	NE-CZ-NH2	-22.07	109.26	120.30
1	D	205	ARG	CD-NE-CZ	13.99	143.18	123.60
1	A	31	ARG	NE-CZ-NH2	10.39	125.50	120.30
1	E	41	ARG	CD-NE-CZ	9.63	137.09	123.60
1	E	31	ARG	NE-CZ-NH2	9.35	124.97	120.30
1	A	41	ARG	NE-CZ-NH1	9.31	124.96	120.30
1	C	31	ARG	NE-CZ-NH2	9.28	124.94	120.30
1	E	205	ARG	NE-CZ-NH1	-9.15	115.73	120.30
1	B	31	ARG	NH1-CZ-NH2	-9.03	109.47	119.40
1	B	205	ARG	NE-CZ-NH1	-8.84	115.88	120.30
1	C	205	ARG	NE-CZ-NH1	-8.76	115.92	120.30
1	E	41	ARG	NE-CZ-NH1	-8.53	116.04	120.30
1	D	41	ARG	NE-CZ-NH1	8.52	124.56	120.30
1	C	41	ARG	NE-CZ-NH1	8.42	124.51	120.30
1	B	41	ARG	NE-CZ-NH1	8.25	124.43	120.30
1	A	205	ARG	NE-CZ-NH1	-8.19	116.20	120.30
1	D	31	ARG	NE-CZ-NH2	8.16	124.38	120.30
1	C	205	ARG	NE-CZ-NH2	7.97	124.29	120.30
1	B	31	ARG	CG-CD-NE	-7.64	95.76	111.80
1	E	205	ARG	NE-CZ-NH2	7.63	124.12	120.30
1	A	31	ARG	NE-CZ-NH1	-7.57	116.51	120.30
1	A	205	ARG	NE-CZ-NH2	7.42	124.01	120.30
1	D	41	ARG	NE-CZ-NH2	-7.36	116.62	120.30
1	B	205	ARG	NE-CZ-NH2	7.18	123.89	120.30
1	A	41	ARG	NE-CZ-NH2	-6.83	116.89	120.30
1	C	41	ARG	NE-CZ-NH2	-6.70	116.95	120.30
1	B	41	ARG	NE-CZ-NH2	-6.63	116.99	120.30
1	C	31	ARG	NE-CZ-NH1	-6.29	117.16	120.30
1	E	298	ARG	NE-CZ-NH2	6.15	123.38	120.30
1	F	344	ASP	CB-CG-OD1	6.10	123.79	118.30
1	E	31	ARG	NE-CZ-NH1	-5.86	117.37	120.30
1	E	141[A]	ARG	NE-CZ-NH1	-5.55	117.53	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	141[B]	ARG	NE-CZ-NH1	-5.55	117.53	120.30
1	A	308	ASP	CB-CG-OD1	5.39	123.15	118.30
1	F	344	ASP	CB-CG-OD2	-5.28	113.55	118.30
1	A	305	ASP	CB-CG-OD1	5.28	123.05	118.30
1	F	440	ASP	CB-CG-OD1	5.20	122.98	118.30
1	D	31	ARG	NE-CZ-NH1	-5.15	117.72	120.30
1	C	344	ASP	CB-CG-OD1	5.10	122.89	118.30
1	D	117	VAL	CG1-CB-CG2	-5.03	102.85	110.90

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	31	ARG	Sidechain
1	D	205	ARG	Sidechain
1	E	41	ARG	Sidechain

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	3591	0	3566	13	0
1	B	3586	0	3562	17	1
1	C	3629	0	3611	12	1
1	D	3622	0	3613	17	0
1	E	3604	0	3576	14	0
1	F	3624	0	3595	42	0
2	A	13	0	18	4	0
2	C	26	0	36	4	0
2	D	13	0	18	2	0
2	F	13	0	18	5	0
3	A	4	0	6	0	0
3	B	16	0	24	0	0
3	C	4	0	6	0	0
3	D	20	0	30	0	0
3	E	12	0	18	0	0
3	F	12	0	18	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	211	0	0	6	0
4	B	171	0	0	4	0
4	C	215	0	0	2	0
4	D	202	0	0	4	0
4	E	174	0	0	3	0
4	F	215	0	0	3	0
All	All	22977	0	21715	123	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 (123) 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)
2:D:467:PG4:H72	2:D:467:PG4:O3	1.65	0.93
1:B:326:VAL:HG11	1:B:362:ALA:HB2	1.53	0.90
1:B:326:VAL:CG1	1:B:362:ALA:HB2	2.10	0.81
1:F:326:VAL:CG1	1:F:360:GLU:HB3	2.11	0.80
1:B:374:GLU:O	1:B:378:VAL:HG23	1.81	0.80
1:F:326:VAL:HG13	1:F:360:GLU:HB3	1.68	0.76
2:D:467:PG4:C7	2:D:467:PG4:O3	2.34	0.75
1:F:277:PHE:HE2	2:F:467:PG4:H32	1.53	0.74
2:A:467:PG4:C7	2:A:467:PG4:O3	2.36	0.72
1:F:117:VAL:HG21	1:F:147:ASN:HB3	1.74	0.69
1:F:104:ALA:HB3	1:F:133:PRO:HB3	1.75	0.69
1:D:313:ARG:HD2	4:D:1180:HOH:O	1.94	0.68
1:F:326:VAL:CG1	1:F:360:GLU:CB	2.71	0.67
1:F:326:VAL:HG13	1:F:360:GLU:CB	2.24	0.67
1:F:163:LEU:HD21	1:F:172:LEU:HD21	1.81	0.62
1:A:313:ARG:HD2	4:A:992:HOH:O	1.99	0.61
1:B:29:GLU:OE1	1:B:201:HIS:NE2	2.30	0.61
1:F:163:LEU:HD21	1:F:172:LEU:HD11	1.82	0.60
1:C:313:ARG:HD2	4:C:1179:HOH:O	1.99	0.60
1:A:458:GLN:NE2	4:A:479:HOH:O	2.32	0.60
1:C:441:GLY:HA2	1:C:462:ILE:HD12	1.83	0.60
1:D:225:ALA:HB1	1:D:300:TRP:CZ3	2.37	0.60
1:F:132:VAL:HB	1:F:133:PRO:CD	2.33	0.58
1:D:19:THR:HG21	1:D:129:LYS:HE2	1.86	0.58
1:D:102:ARG:NH1	1:D:289:GLU:OE1	2.37	0.58
1:D:330:LYS:HE3	1:D:365:HIS:CE1	2.40	0.57
1:F:273:GLY:H	2:F:467:PG4:H22	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:182:GLY:HA2	1:F:213:THR:HG22	1.88	0.55
1:B:155:GLN:OE1	4:B:683:HOH:O	2.18	0.55
1:F:27:CYS:HB3	1:F:30:ILE:HD12	1.87	0.54
1:B:135:ARG:HA	4:B:472:HOH:O	2.07	0.54
1:F:373:ARG:NH1	4:F:484:HOH:O	2.40	0.54
1:F:211:THR:HG23	1:F:212:ASN:O	2.07	0.54
1:F:26:MET:CE	1:F:173:LYS:HG2	2.38	0.53
1:D:313:ARG:CD	4:D:1180:HOH:O	2.54	0.53
1:A:31:ARG:HD3	4:A:491:HOH:O	2.07	0.53
1:A:29:GLU:OE1	1:A:201:HIS:NE2	2.33	0.52
1:A:313:ARG:NH1	4:A:992:HOH:O	1.94	0.52
1:C:277:PHE:HE2	2:C:467:PG4:H22	1.75	0.52
1:F:135:ARG:HA	1:F:138:GLU:OE2	2.10	0.52
1:F:19:THR:HG21	1:F:129:LYS:HE3	1.91	0.52
1:F:277:PHE:CE2	2:F:467:PG4:H32	2.39	0.52
1:C:0:SER:O	1:C:1:MET:HB3	2.10	0.51
1:F:39:GLU:HG3	1:F:73:THR:HG21	1.92	0.51
1:C:281:VAL:O	1:C:285:VAL:HG23	2.11	0.51
1:D:325:THR:O	1:D:325:THR:HG23	2.11	0.50
1:F:84:LEU:C	1:F:84:LEU:HD13	2.33	0.49
1:C:443:ARG:HD3	1:C:462:ILE:O	2.12	0.49
1:E:77:ASP:HB3	4:E:1026:HOH:O	2.13	0.49
1:E:14:TYR:OH	1:E:165:GLU:HA	2.12	0.49
1:F:224:ASN:ND2	1:F:280:ASP:OD2	2.47	0.48
1:B:325:THR:O	1:B:325:THR:HG23	2.13	0.48
1:C:330:LYS:HE3	1:C:365:HIS:CE1	2.48	0.48
4:C:528:HOH:O	1:E:142:ARG:CD	2.62	0.48
1:F:382:HIS:ND1	1:F:388:ASP:OD1	2.45	0.48
1:E:330:LYS:HE3	1:E:365:HIS:CE1	2.49	0.47
2:C:467:PG4:H42	2:C:467:PG4:H11	1.96	0.47
1:E:335:GLY:HA3	1:E:415:THR:HB	1.95	0.47
1:D:102:ARG:HD3	4:D:511:HOH:O	2.15	0.47
1:B:374:GLU:HB2	4:B:1154:HOH:O	2.14	0.47
1:F:132:VAL:CB	1:F:133:PRO:CD	2.93	0.47
1:E:458:GLN:NE2	4:E:498:HOH:O	2.35	0.47
1:D:134:VAL:O	1:D:135:ARG:HB2	2.15	0.46
2:C:467:PG4:H42	2:C:467:PG4:C1	2.46	0.46
1:E:389:ASP:O	1:E:393:ARG:HB2	2.14	0.46
1:B:330:LYS:HE3	1:B:365:HIS:CE1	2.50	0.46
1:B:141[A]:ARG:NH1	4:B:533:HOH:O	2.48	0.46
1:A:330:LYS:HE3	1:A:365:HIS:CE1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:LYS:HE3	1:A:365:HIS:CG	2.52	0.45
1:E:246:ALA:CB	1:F:218:LEU:HD23	2.45	0.45
2:C:467:PG4:C4	2:C:467:PG4:C1	2.95	0.45
1:C:189:GLN:O	1:C:193:GLN:HG2	2.16	0.45
1:F:102:ARG:HD2	4:F:884:HOH:O	2.16	0.45
2:A:467:PG4:O3	2:A:467:PG4:H71	2.17	0.45
1:B:326:VAL:CG1	1:B:362:ALA:CB	2.89	0.45
1:D:3:GLU:OE2	1:D:5:LYS:HE3	2.17	0.45
1:F:464:LYS:NZ	4:F:586:HOH:O	2.50	0.45
1:F:26:MET:HE2	1:F:173:LYS:HG2	1.99	0.44
1:F:390:GLN:HG2	1:F:394:LEU:HD12	1.98	0.44
1:F:84:LEU:HD13	1:F:85:VAL:N	2.33	0.44
1:B:19:THR:HG21	1:B:129:LYS:HE2	1.99	0.44
1:E:186:PRO:O	1:E:190[A]:ARG:HG3	2.17	0.44
1:D:76:ASP:OD1	1:D:115:ARG:NE	2.46	0.44
1:A:373:ARG:HB2	1:A:397:ILE:HD12	1.99	0.44
1:C:1:MET:CG	1:C:1:MET:O	2.66	0.44
1:E:309:TYR:CZ	1:E:313:ARG:HD3	2.53	0.44
1:C:134:VAL:O	1:C:135:ARG:HB2	2.18	0.44
1:F:106:LEU:O	1:F:107:LYS:C	2.56	0.44
1:A:434:LYS:HA	1:A:435:PRO:C	2.37	0.43
1:A:330:LYS:HE3	1:A:365:HIS:CD2	2.53	0.43
1:B:79:ILE:O	1:B:124:LYS:HE3	2.17	0.43
1:B:326:VAL:HG12	1:B:326:VAL:O	2.18	0.43
1:F:26:MET:HE3	1:F:173:LYS:HG2	2.00	0.43
1:F:132:VAL:HB	1:F:133:PRO:HD2	2.00	0.43
1:A:405:CYS:HB3	1:A:428:ILE:HG23	2.01	0.43
1:B:368:ASP:OD1	1:B:371:VAL:HG23	2.19	0.43
1:F:132:VAL:HG11	1:F:279:LYS:HD3	2.00	0.43
1:E:267:LYS:HB3	4:E:912:HOH:O	2.19	0.43
1:F:273:GLY:H	2:F:467:PG4:C2	2.31	0.42
1:C:19:THR:HG21	1:C:129:LYS:HE2	2.01	0.42
1:F:131:ALA:O	1:F:132:VAL:O	2.37	0.42
1:D:313:ARG:CG	4:D:1180:HOH:O	2.67	0.42
1:A:332:ALA:HB2	1:A:404:ALA:O	2.19	0.42
1:F:134:VAL:O	1:F:135:ARG:HB2	2.20	0.42
1:F:326:VAL:HG13	1:F:360:GLU:HB2	2.00	0.41
1:F:84:LEU:C	1:F:84:LEU:CD1	2.89	0.41
2:A:467:PG4:H22	4:A:761:HOH:O	2.21	0.41
2:A:467:PG4:O3	2:A:467:PG4:H72	2.20	0.41
1:F:373:ARG:HB2	1:F:397:ILE:HG21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:326:VAL:HG12	1:B:362:ALA:CA	2.51	0.41
1:A:393:ARG:NH1	4:A:808:HOH:O	2.54	0.41
1:D:6:LYS:HD3	1:D:81:GLU:HB3	2.03	0.41
1:F:273:GLY:N	2:F:467:PG4:H22	2.36	0.40
1:F:132:VAL:O	1:F:133:PRO:C	2.60	0.40
1:D:225:ALA:HB1	1:D:300:TRP:CE3	2.56	0.40
1:D:188:GLY:O	1:D:192:VAL:HG23	2.20	0.40
1:D:326:VAL:O	1:D:326:VAL:HG12	2.21	0.40
1:C:434:LYS:HA	1:C:435:PRO:C	2.41	0.40
1:D:269:SER:C	1:D:462:ILE:HD13	2.41	0.40
1:E:325:THR:HG23	1:E:325:THR:O	2.20	0.40
1:E:443:ARG:HD3	1:E:462:ILE:O	2.21	0.40
1:E:443:ARG:HD2	1:E:461:THR:OG1	2.22	0.40
1:B:189:GLN:O	1:B:193:GLN:HG2	2.21	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:B:40:SER:OG	1:C:77:ASP:OD1[1_455]	1.82	0.38

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	460/467 (98%)	449 (98%)	11 (2%)	0	100	100
1	B	460/467 (98%)	447 (97%)	13 (3%)	0	100	100
1	C	466/467 (100%)	453 (97%)	13 (3%)	0	100	100
1	D	465/467 (100%)	455 (98%)	10 (2%)	0	100	100
1	E	466/467 (100%)	454 (97%)	12 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	465/467 (100%)	446 (96%)	18 (4%)	1 (0%)	52	69
All	All	2782/2802 (99%)	2704 (97%)	77 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	132	VAL

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	385/400 (96%)	382 (99%)	3 (1%)	86	94
1	B	386/400 (96%)	383 (99%)	3 (1%)	86	94
1	C	391/400 (98%)	388 (99%)	3 (1%)	86	94
1	D	391/400 (98%)	389 (100%)	2 (0%)	92	97
1	E	385/400 (96%)	381 (99%)	4 (1%)	82	93
1	F	391/400 (98%)	387 (99%)	4 (1%)	82	93
All	All	2329/2400 (97%)	2310 (99%)	19 (1%)	86	94

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	GLU
1	A	177	ARG
1	A	449	HIS
1	B	161	GLU
1	B	177	ARG
1	B	374	GLU
1	C	161	GLU
1	C	177	ARG
1	C	184	GLU
1	D	161	GLU

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Mol	Chain	Res	Type
1	D	177	ARG
1	E	117	VAL
1	E	161	GLU
1	E	177	ARG
1	E	459	ILE
1	F	84	LEU
1	F	129	LYS
1	F	177	ARG
1	F	206	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	121	ASN
1	B	302	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

22 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	PG4	A	467	-	12,12,12	0.53	0	11,11,11	0.58	0
3	EDO	A	468	-	3,3,3	0.63	0	2,2,2	0.16	0
3	EDO	B	467	-	3,3,3	0.60	0	2,2,2	0.23	0
3	EDO	B	468	-	3,3,3	0.53	0	2,2,2	0.35	0
3	EDO	B	469	-	3,3,3	0.49	0	2,2,2	0.60	0
3	EDO	B	470	-	3,3,3	0.57	0	2,2,2	0.19	0
2	PG4	C	467	-	12,12,12	0.61	0	11,11,11	0.85	0
2	PG4	C	468	-	12,12,12	0.49	0	11,11,11	0.46	0
3	EDO	C	469	-	3,3,3	0.63	0	2,2,2	0.18	0
2	PG4	D	467	-	12,12,12	0.83	0	11,11,11	1.05	0
3	EDO	D	468	-	3,3,3	0.71	0	2,2,2	0.23	0
3	EDO	D	469	-	3,3,3	0.55	0	2,2,2	0.40	0
3	EDO	D	470	-	3,3,3	0.54	0	2,2,2	0.74	0
3	EDO	D	471	-	3,3,3	0.61	0	2,2,2	0.18	0
3	EDO	D	472	-	3,3,3	0.54	0	2,2,2	0.30	0
3	EDO	E	467	-	3,3,3	0.69	0	2,2,2	0.03	0
3	EDO	E	468	-	3,3,3	0.70	0	2,2,2	0.13	0
3	EDO	E	469	-	3,3,3	0.58	0	2,2,2	0.21	0
2	PG4	F	467	-	12,12,12	0.67	0	11,11,11	0.85	0
3	EDO	F	468	-	3,3,3	0.53	0	2,2,2	0.24	0
3	EDO	F	469	-	3,3,3	0.66	0	2,2,2	0.24	0
3	EDO	F	470	-	3,3,3	0.60	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
2	PG4	A	467	-	-	0/10/10/10	0/0/0/0
3	EDO	A	468	-	-	0/1/1/1	0/0/0/0
3	EDO	B	467	-	-	0/1/1/1	0/0/0/0
3	EDO	B	468	-	-	0/1/1/1	0/0/0/0
3	EDO	B	469	-	-	0/1/1/1	0/0/0/0
3	EDO	B	470	-	-	0/1/1/1	0/0/0/0
2	PG4	C	467	-	-	0/10/10/10	0/0/0/0
2	PG4	C	468	-	-	0/10/10/10	0/0/0/0
3	EDO	C	469	-	-	0/1/1/1	0/0/0/0
2	PG4	D	467	-	-	0/10/10/10	0/0/0/0
3	EDO	D	468	-	-	0/1/1/1	0/0/0/0
3	EDO	D	469	-	-	0/1/1/1	0/0/0/0
3	EDO	D	470	-	-	0/1/1/1	0/0/0/0
3	EDO	D	471	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	D	472	-	-	0/1/1/1	0/0/0/0
3	EDO	E	467	-	-	0/1/1/1	0/0/0/0
3	EDO	E	468	-	-	0/1/1/1	0/0/0/0
3	EDO	E	469	-	-	0/1/1/1	0/0/0/0
2	PG4	F	467	-	-	0/10/10/10	0/0/0/0
3	EDO	F	468	-	-	0/1/1/1	0/0/0/0
3	EDO	F	469	-	-	0/1/1/1	0/0/0/0
3	EDO	F	470	-	-	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.

4 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	467	PG4	4	0
2	C	467	PG4	4	0
2	D	467	PG4	2	0
2	F	467	PG4	5	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	464/467 (99%)	-0.62	2 (0%) 93 93	2, 13, 32, 52	0
1	B	462/467 (98%)	-0.44	4 (0%) 85 85	3, 13, 32, 58	0
1	C	467/467 (100%)	-0.59	3 (0%) 90 90	3, 13, 32, 68	0
1	D	467/467 (100%)	-0.40	5 (1%) 82 82	3, 13, 32, 65	0
1	E	466/467 (99%)	-0.40	3 (0%) 90 90	2, 13, 32, 64	0
1	F	467/467 (100%)	-0.61	5 (1%) 82 82	3, 13, 32, 64	0
All	All	2793/2802 (99%)	-0.51	22 (0%) 87 87	2, 13, 32, 68	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	0	SER	6.0
1	A	1	MET	5.0
1	B	389	ASP	4.8
1	F	0	SER	4.3
1	F	384	GLY	4.1
1	C	1	MET	4.0
1	E	384	GLY	4.0
1	D	386	SER	3.7
1	C	384	GLY	3.5
1	D	0	SER	3.2
1	B	0	SER	3.2
1	E	385	VAL	3.1
1	A	2	PHE	3.0
1	E	386	SER	2.8
1	F	131	ALA	2.6
1	D	383	PRO	2.5
1	F	132	VAL	2.4
1	F	134	VAL	2.4
1	D	384	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	391	VAL	2.2
1	B	164	ALA	2.2
1	D	164	ALA	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	EDO	D	470	4/4	0.59	0.34	12.30	43,56,58,63	0
2	PG4	F	467	13/13	0.83	0.26	5.98	30,47,61,67	0
3	EDO	F	469	4/4	0.83	0.26	5.54	39,45,49,51	0
2	PG4	C	467	13/13	0.89	0.19	5.49	26,42,49,49	0
3	EDO	F	468	4/4	0.87	0.24	5.34	36,47,55,57	0
3	EDO	B	469	4/4	0.86	0.24	5.28	29,51,56,60	0
3	EDO	D	468	4/4	0.92	0.16	4.55	38,39,41,42	0
2	PG4	A	467	13/13	0.95	0.15	3.99	17,36,51,54	0
2	PG4	D	467	13/13	0.85	0.21	3.57	27,44,51,55	0
3	EDO	B	468	4/4	0.91	0.21	1.77	52,59,63,64	0
2	PG4	C	468	13/13	0.91	0.17	1.56	23,49,59,64	0
3	EDO	D	471	4/4	0.88	0.14	0.74	39,49,49,53	0
3	EDO	A	468	4/4	0.90	0.13	0.68	32,39,44,48	0
3	EDO	E	468	4/4	0.76	0.18	0.54	49,54,55,61	0
3	EDO	E	467	4/4	0.91	0.28	-	32,37,37,38	0
3	EDO	B	467	4/4	0.75	0.18	-	58,59,59,65	0
3	EDO	B	470	4/4	0.88	0.27	-	38,54,55,57	0
3	EDO	C	469	4/4	0.72	0.20	-	55,63,65,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	EDO	D	469	4/4	0.76	0.27	-	60,63,69,69	0
3	EDO	E	469	4/4	0.88	0.15	-	49,50,54,56	0
3	EDO	D	472	4/4	0.88	0.15	-	62,64,66,70	0
3	EDO	F	470	4/4	0.82	0.16	-	52,53,55,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.