



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2018 – 11:27 pm GMT

PDB ID : 4F9F
Title : Crystal Structure of VldE, the pseudo-glycosyltransferase, in complex with GDP and Trehalose
Authors : Cavalier, M.C.; Yim, Y.-S.; Asamizu, S.; Neau, D.; Almabruk, K.H.; Mahmud, T.; Lee, Y.-H.
Deposited on : 2012-05-18
Resolution : 2.81 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

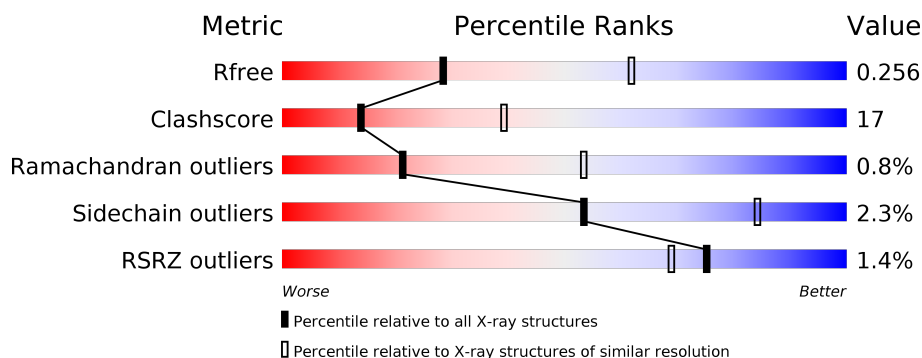
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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}	111664	2792 (2.80-2.80)
Clashscore	122126	3209 (2.80-2.80)
Ramachandran outliers	120053	3158 (2.80-2.80)
Sidechain outliers	120020	3160 (2.80-2.80)
RSRZ outliers	108989	2726 (2.80-2.80)

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	497	<div> <div>69%</div> <div>24%</div> <div>• 5%</div> </div>
1	B	497	<div> <div>%</div> <div>70%</div> <div>21%</div> <div>• 7%</div> </div>
1	C	497	<div> <div>2%</div> <div>67%</div> <div>24%</div> <div>• 7%</div> </div>
1	D	497	<div> <div>%</div> <div>71%</div> <div>22%</div> <div>• 7%</div> </div>
1	E	497	<div> <div>%</div> <div>68%</div> <div>23%</div> <div>• 7%</div> </div>
1	F	497	<div> <div>2%</div> <div>69%</div> <div>21%</div> <div>• 7%</div> </div>

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	GDP	A	501	-	-	X	-
3	TRE	A	502	-	-	X	-

2 Entry composition [i](#)

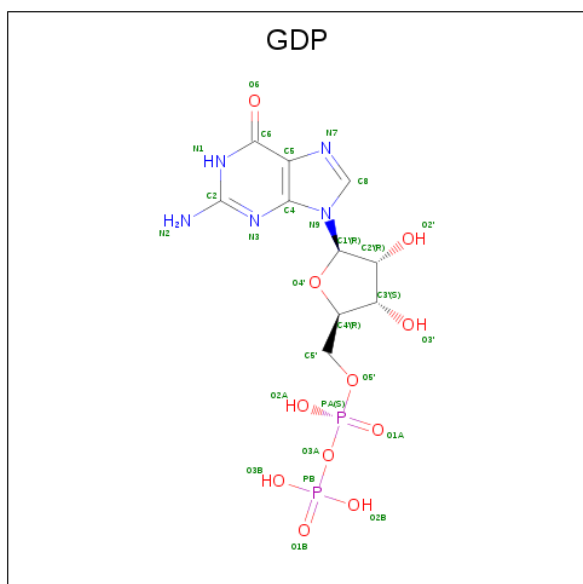
There are 4 unique types of molecules in this entry. The entry contains 22212 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 VldE.

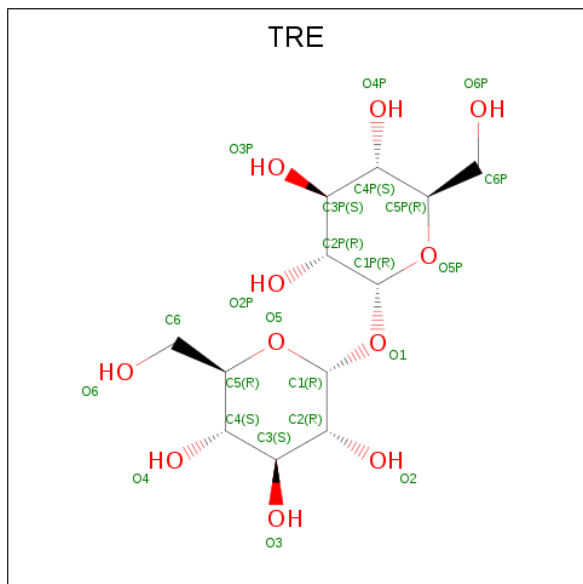
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	470	Total	C	N	O	S	0	0	0
			3674	2301	675	686	12			
1	B	463	Total	C	N	O	S	0	0	0
			3636	2278	671	675	12			
1	C	460	Total	C	N	O	S	0	0	0
			3605	2260	663	670	12			
1	D	464	Total	C	N	O	S	0	0	0
			3638	2279	672	675	12			
1	E	463	Total	C	N	O	S	0	0	0
			3633	2276	671	674	12			
1	F	460	Total	C	N	O	S	0	0	0
			3610	2261	667	670	12			

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	C	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	D	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	E	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	F	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 3 is TREHALOSE (three-letter code: TRE) (formula: $C_{12}H_{22}O_{11}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			23	12	11		
3	D	1	Total	C	O	0	0
			23	12	11		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	36	Total	O	0	0
			36	36		
4	B	22	Total	O	0	0
			22	22		

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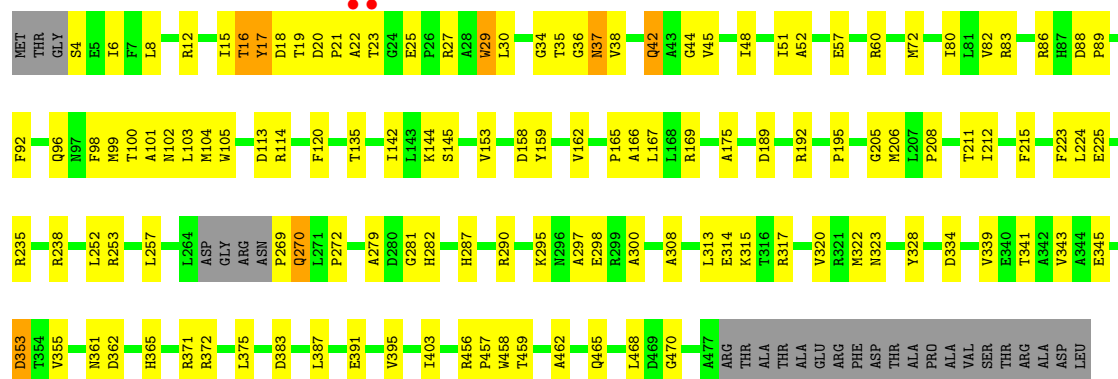
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	31	Total 31	O 31	0	0
4	D	44	Total 44	O 44	0	0
4	E	41	Total 41	O 41	0	0
4	F	28	Total 28	O 28	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

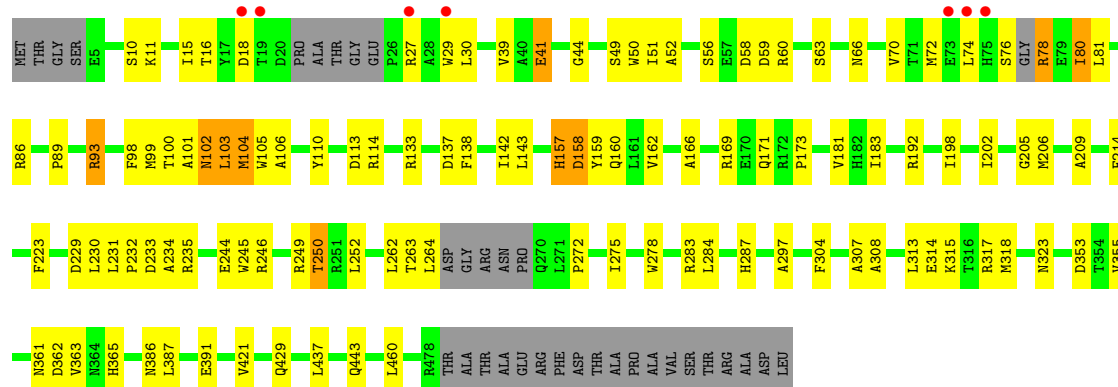
• Molecule 1: VldE

Chain A: 



• Molecule 1: VldE

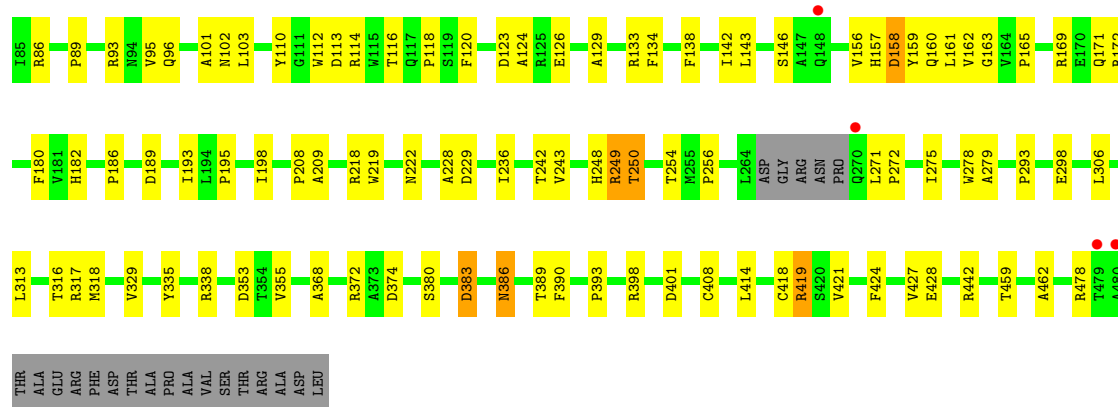
Chain B: 



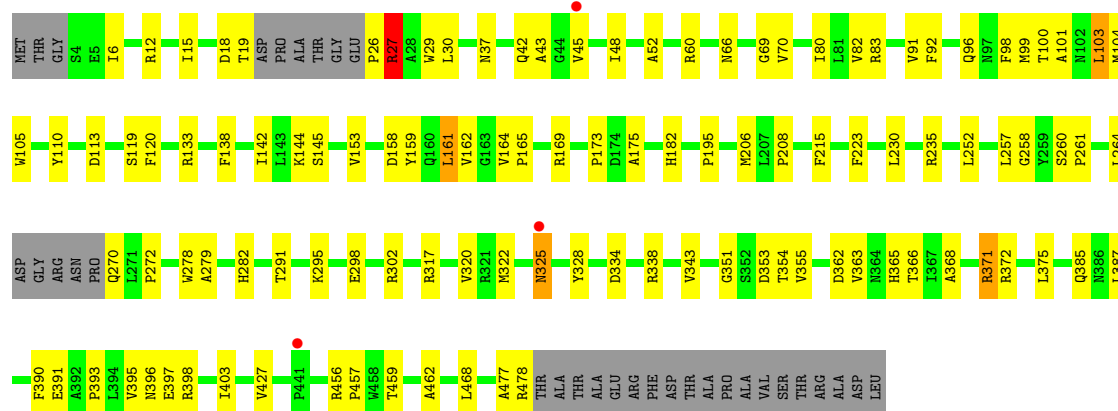
• Molecule 1: VldE

Chain C: 

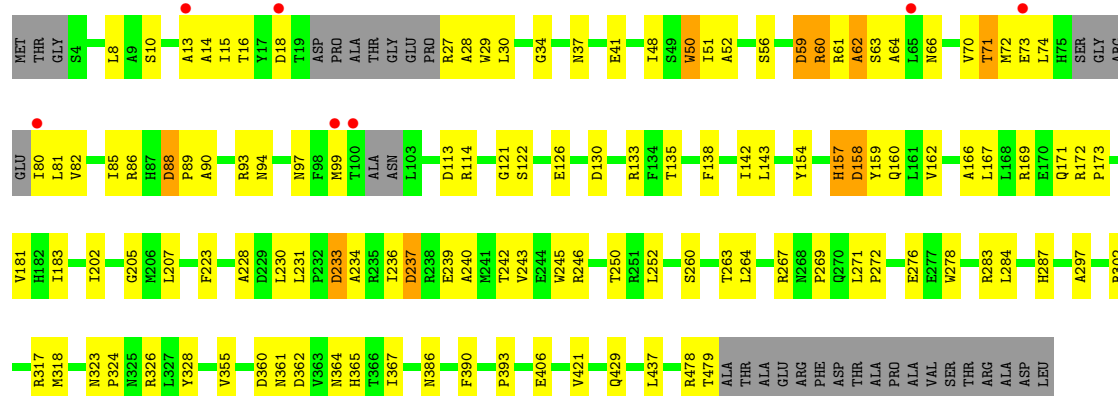




• Molecule 1: VldE

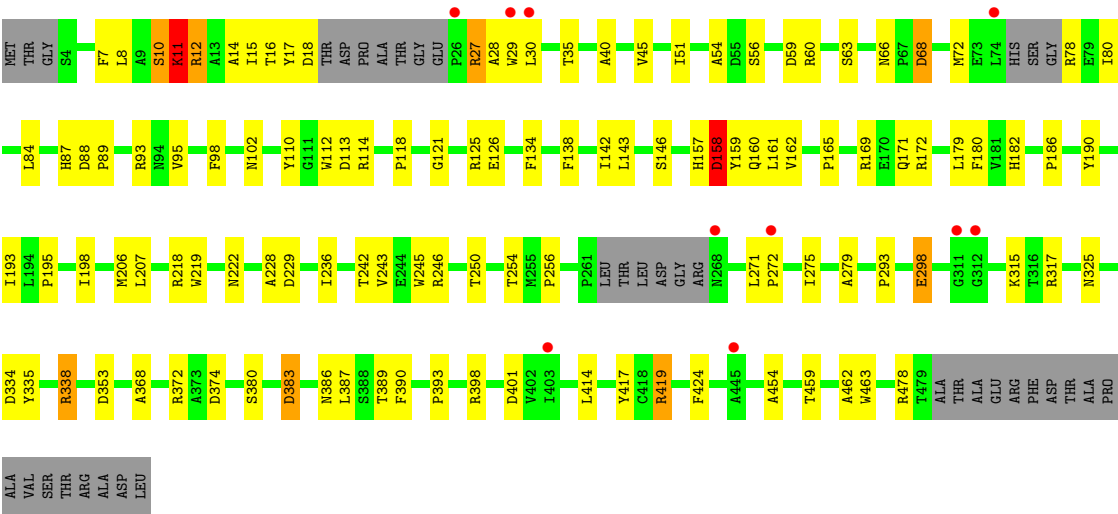


• Molecule 1: VldE



• Molecule 1: VldE





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	320.82Å 122.82Å 93.86Å 90.00° 91.62° 90.00°	Depositor
Resolution (Å)	49.15 – 2.81 49.15 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.15-2.81) 92.4 (49.15-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.211 , 0.258 0.208 , 0.256	Depositor DCC
R_{free} test set	4425 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	22212	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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: GDP, TRE

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	0.47	0/3760	0.61	0/5125
1	B	0.47	0/3718	0.59	0/5062
1	C	0.44	0/3685	0.60	0/5018
1	D	0.47	0/3721	0.61	2/5067 (0.0%)
1	E	0.47	0/3715	0.60	0/5059
1	F	0.44	0/3692	0.60	2/5027 (0.0%)
All	All	0.46	0/22291	0.60	4/30358 (0.0%)

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	C	0	1
1	E	0	1
1	F	0	1
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	43	ALA	CB-CA-C	-6.34	100.59	110.10
1	F	11	LYS	N-CA-C	5.29	125.30	111.00
1	D	43	ALA	N-CA-C	5.17	124.95	111.00
1	F	419	ARG	O-C-N	-5.02	114.67	122.70

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	157	HIS	Peptide
1	C	13	ALA	Peptide
1	E	157	HIS	Peptide
1	F	338	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	3674	0	3565	131	0
1	B	3636	0	3533	130	1
1	C	3605	0	3506	131	1
1	D	3638	0	3538	100	2
1	E	3633	0	3531	125	0
1	F	3610	0	3506	126	0
2	A	28	0	12	15	0
2	B	28	0	12	3	0
2	C	28	0	12	0	0
2	D	28	0	12	7	0
2	E	28	0	12	0	0
2	F	28	0	12	2	0
3	A	23	0	22	10	0
3	D	23	0	22	4	0
4	A	36	0	0	5	1
4	B	22	0	0	4	0
4	C	31	0	0	1	1
4	D	44	0	0	4	0
4	E	41	0	0	4	0
4	F	28	0	0	1	0
All	All	22212	0	21295	732	3

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 17.

The worst 5 of 732 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:74:LEU:HD11	1:C:80:ILE:CD1	1.53	1.37
2:A:501:GDP:O3B	3:A:502:TRE:HC3	1.12	1.28
1:A:317:ARG:NH1	1:A:353:ASP:O	1.68	1.25
1:A:169:ARG:NH2	1:A:175:ALA:O	1.65	1.24
1:A:269:PRO:O	1:A:270:GLN:HG2	1.37	1.24

All (3) 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:ARG:NH2	1:D:397:GLU:OE1[4_445]	1.75	0.45
1:D:264:LEU:C	4:C:721:HOH:O[4_455]	2.09	0.11
1:B:443:GLN:OE1	4:A:616:HOH:O[2_555]	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	466/497 (94%)	438 (94%)	22 (5%)	6 (1%)	13	40
1	B	455/497 (92%)	426 (94%)	26 (6%)	3 (1%)	24	57
1	C	452/497 (91%)	428 (95%)	21 (5%)	3 (1%)	24	57
1	D	458/497 (92%)	435 (95%)	20 (4%)	3 (1%)	24	57
1	E	455/497 (92%)	427 (94%)	25 (6%)	3 (1%)	24	57
1	F	452/497 (91%)	426 (94%)	23 (5%)	3 (1%)	24	57
All	All	2738/2982 (92%)	2580 (94%)	137 (5%)	21 (1%)	21	53

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	158	ASP
1	E	158	ASP

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Mol	Chain	Res	Type
1	A	458	TRP
1	C	158	ASP
1	E	386	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	375/395 (95%)	368 (98%)	7 (2%)	60	87
1	B	371/395 (94%)	362 (98%)	9 (2%)	52	83
1	C	367/395 (93%)	360 (98%)	7 (2%)	60	87
1	D	371/395 (94%)	365 (98%)	6 (2%)	65	90
1	E	371/395 (94%)	360 (97%)	11 (3%)	44	78
1	F	368/395 (93%)	357 (97%)	11 (3%)	44	78
All	All	2223/2370 (94%)	2172 (98%)	51 (2%)	53	84

5 of 51 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	27	ARG
1	D	371	ARG
1	F	298	GLU
1	D	161	LEU
1	E	27	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	109	ASN
1	D	182	HIS
1	E	96	GLN
1	C	157	HIS
1	E	75	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 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	GDP	A	501	-	25,30,30	1.23	2 (8%)	27,47,47	2.11	7 (25%)
3	TRE	A	502	-	24,24,24	1.69	5 (20%)	35,35,35	1.12	3 (8%)
2	GDP	B	600	-	25,30,30	1.22	2 (8%)	27,47,47	2.12	7 (25%)
2	GDP	C	600	-	25,30,30	1.23	2 (8%)	27,47,47	2.11	7 (25%)
2	GDP	D	501	-	25,30,30	1.23	2 (8%)	27,47,47	2.13	7 (25%)
3	TRE	D	502	-	24,24,24	1.72	5 (20%)	35,35,35	1.32	4 (11%)
2	GDP	E	600	-	25,30,30	1.24	2 (8%)	27,47,47	2.12	7 (25%)
2	GDP	F	600	-	25,30,30	1.21	2 (8%)	27,47,47	2.12	7 (25%)

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	GDP	A	501	-	-	0/12/32/32	0/3/3/3
3	TRE	A	502	-	-	0/8/48/48	0/2/2/2
2	GDP	B	600	-	-	0/12/32/32	0/3/3/3
2	GDP	C	600	-	-	0/12/32/32	0/3/3/3
2	GDP	D	501	-	-	0/12/32/32	0/3/3/3
3	TRE	D	502	-	-	0/8/48/48	0/2/2/2
2	GDP	E	600	-	-	0/12/32/32	0/3/3/3
2	GDP	F	600	-	-	0/12/32/32	0/3/3/3

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	502	TRE	C4P-C3P	-3.37	1.43	1.52
3	A	502	TRE	C4P-C3P	-3.34	1.43	1.52
3	A	502	TRE	C3-C2	-2.58	1.45	1.52
3	D	502	TRE	C3-C2	-2.53	1.45	1.52
3	D	502	TRE	C3P-C2P	-2.30	1.46	1.52

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	GDP	C5-C6-N1	-3.86	117.98	123.47
2	B	600	GDP	C5-C6-N1	-3.85	117.99	123.47
2	D	501	GDP	C5-C6-N1	-3.85	118.00	123.47
2	C	600	GDP	C5-C6-N1	-3.84	118.00	123.47
2	F	600	GDP	C5-C6-N1	-3.83	118.03	123.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 31 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	GDP	15	0
3	A	502	TRE	10	0
2	B	600	GDP	3	0
2	D	501	GDP	7	0
3	D	502	TRE	4	0
2	F	600	GDP	2	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 [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	470/497 (94%)	-0.20	2 (0%) 92 90	26, 54, 101, 192	0
1	B	463/497 (93%)	-0.08	7 (1%) 73 67	28, 56, 111, 167	0
1	C	460/497 (92%)	-0.09	9 (1%) 65 56	27, 58, 103, 157	0
1	D	464/497 (93%)	-0.23	3 (0%) 89 86	28, 53, 94, 151	0
1	E	463/497 (93%)	-0.14	7 (1%) 73 67	27, 53, 108, 162	0
1	F	460/497 (92%)	0.04	10 (2%) 62 52	34, 64, 110, 165	0
All	All	2780/2982 (93%)	-0.12	38 (1%) 75 69	26, 57, 106, 192	0

The worst 5 of 38 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	480	ALA	5.3
1	F	268	ASN	4.2
1	A	22	ALA	4.0
1	B	73	GLU	3.5
1	C	4	SER	3.4

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. 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	TRE	D	502	23/23	0.86	0.27	62,62,62,62	23
2	GDP	A	501	28/28	0.87	0.18	62,62,62,62	28
3	TRE	A	502	23/23	0.88	0.24	62,62,62,62	23
2	GDP	F	600	28/28	0.91	0.20	62,62,62,62	0
2	GDP	C	600	28/28	0.91	0.19	62,62,62,62	0
2	GDP	D	501	28/28	0.92	0.18	62,62,62,62	28
2	GDP	B	600	28/28	0.94	0.16	62,62,62,62	0
2	GDP	E	600	28/28	0.94	0.21	62,62,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.