



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 15, 2026 – 11:02 AM UTC

PDB ID : 6DQG / pdb_00006dqg
Title : Human glutamate dehydrogenase, H454Y mutant
Authors : Smith, T.J.
Deposited on : 2018-06-10
Resolution : 2.70 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 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

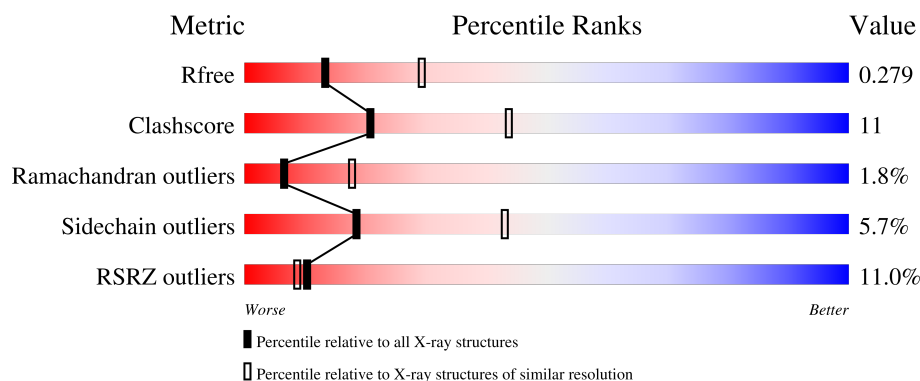
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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.



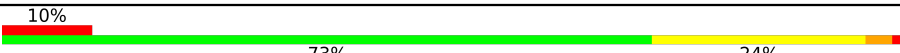
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	496	8% 77% 20% .
1	B	496	13% 72% 24% .
1	C	496	23% 74% 23% ..
1	D	496	10% 72% 24% .
1	E	496	3% 73% 23% .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	496	

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	PO4	A	1001	-	-	X	-
2	PO4	D	603	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 23591 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 Glutamate dehydrogenase 1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	496	Total	C	N	O	S	0	0	0
			3876	2453	677	727	19			
1	B	496	Total	C	N	O	S	0	0	0
			3876	2453	677	727	19			
1	C	496	Total	C	N	O	S	0	0	0
			3876	2453	677	727	19			
1	D	496	Total	C	N	O	S	0	0	0
			3876	2453	677	727	19			
1	E	496	Total	C	N	O	S	0	0	0
			3876	2453	677	727	19			
1	F	496	Total	C	N	O	S	0	0	0
			3876	2453	677	727	19			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	454	TYR	HIS	engineered mutation	UNP P00367
B	454	TYR	HIS	engineered mutation	UNP P00367
C	454	TYR	HIS	engineered mutation	UNP P00367
D	454	TYR	HIS	engineered mutation	UNP P00367
E	454	TYR	HIS	engineered mutation	UNP P00367
F	454	TYR	HIS	engineered mutation	UNP P00367

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	O	P	0	0
			5	4	1		

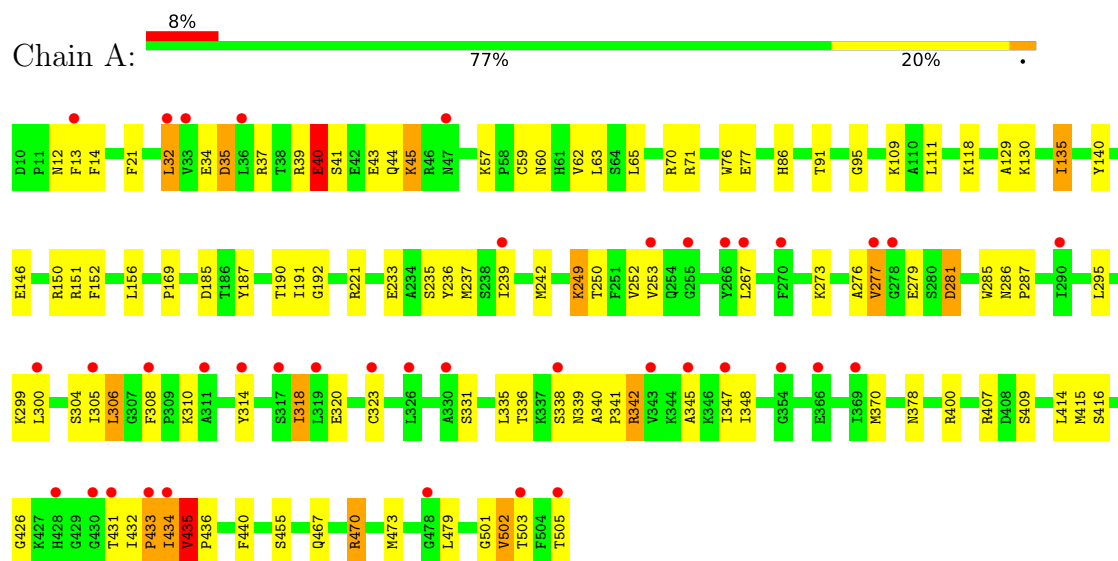
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	31	Total	O	0	0
			31	31		
3	B	51	Total	O	0	0
			51	51		
3	C	46	Total	O	0	0
			46	46		
3	D	49	Total	O	0	0
			49	49		
3	E	47	Total	O	0	0
			47	47		
3	F	36	Total	O	0	0
			36	36		

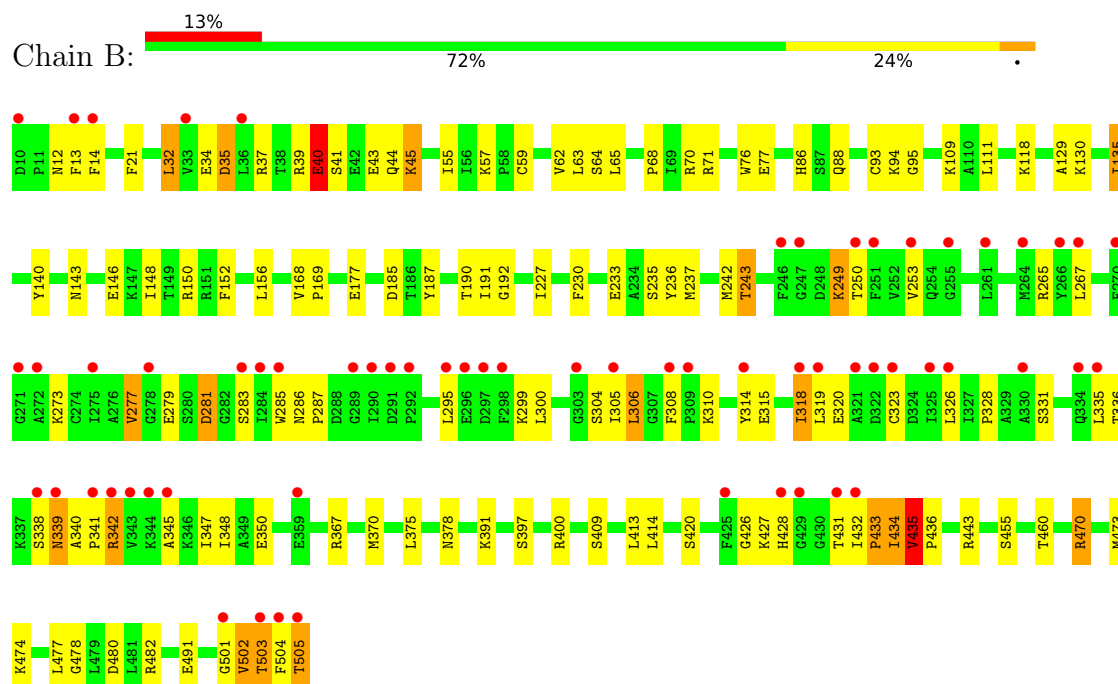
3 Residue-property plots

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.

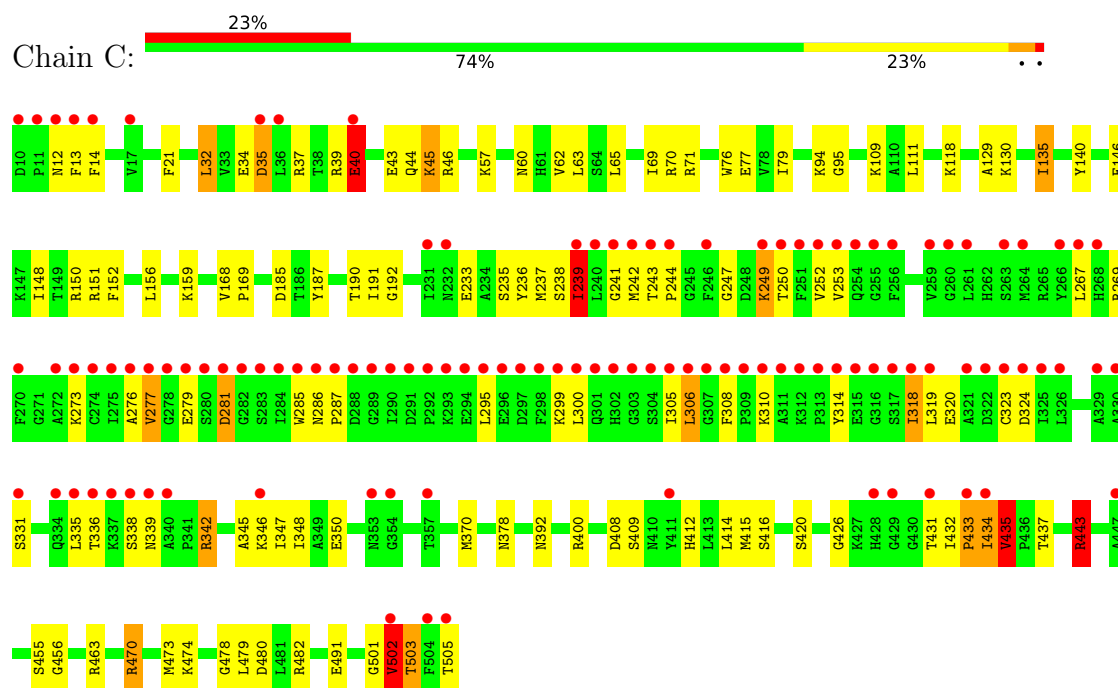
- Molecule 1: Glutamate dehydrogenase 1, mitochondrial



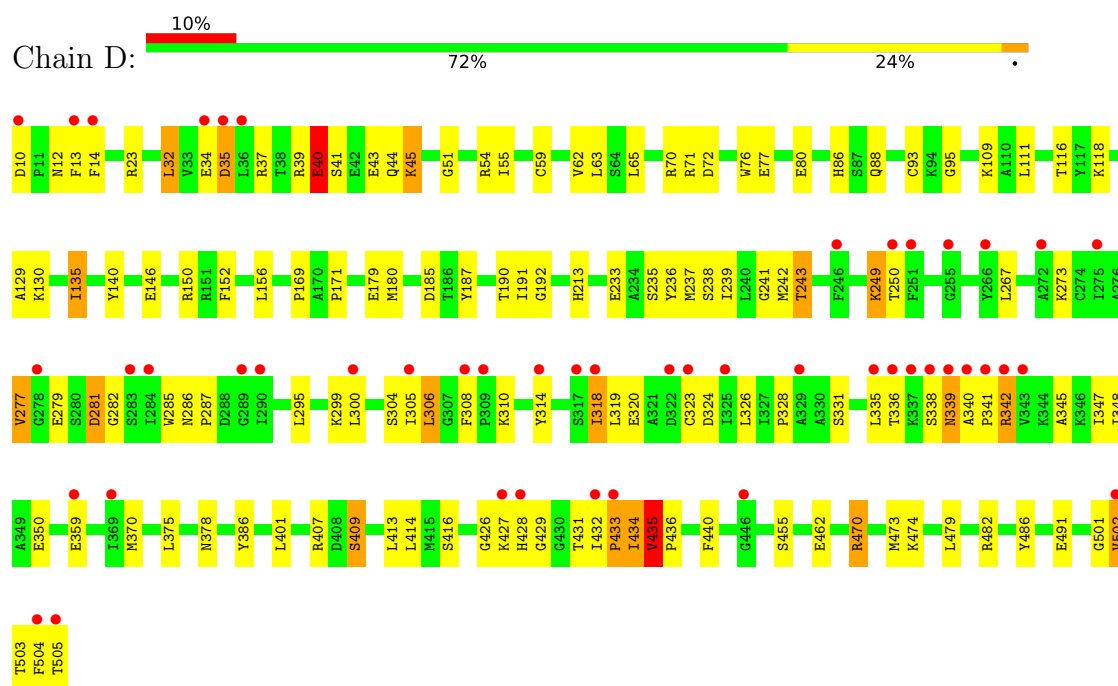
- Molecule 1: Glutamate dehydrogenase 1, mitochondrial



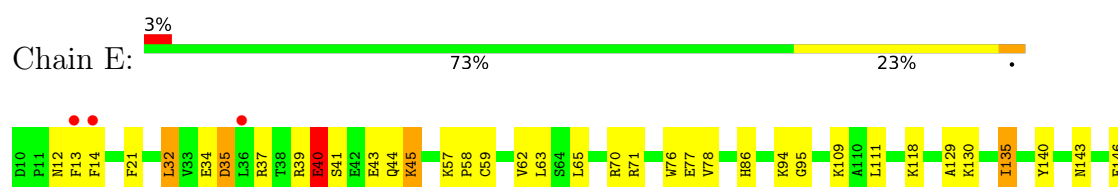
• Molecule 1: Glutamate dehydrogenase 1, mitochondrial

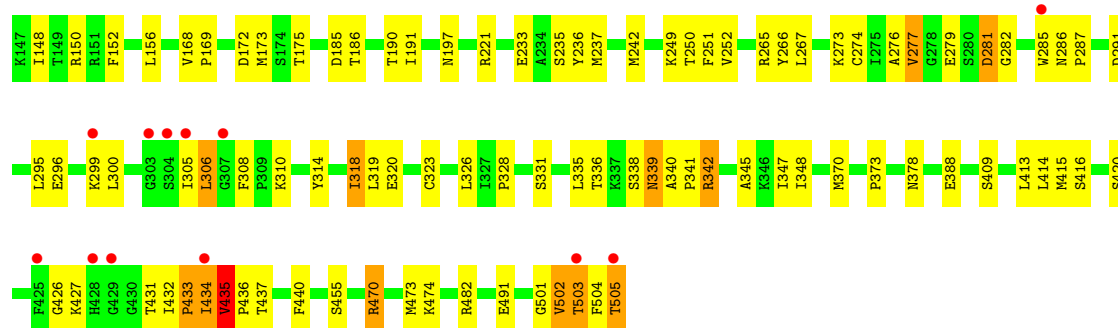


• Molecule 1: Glutamate dehydrogenase 1, mitochondrial

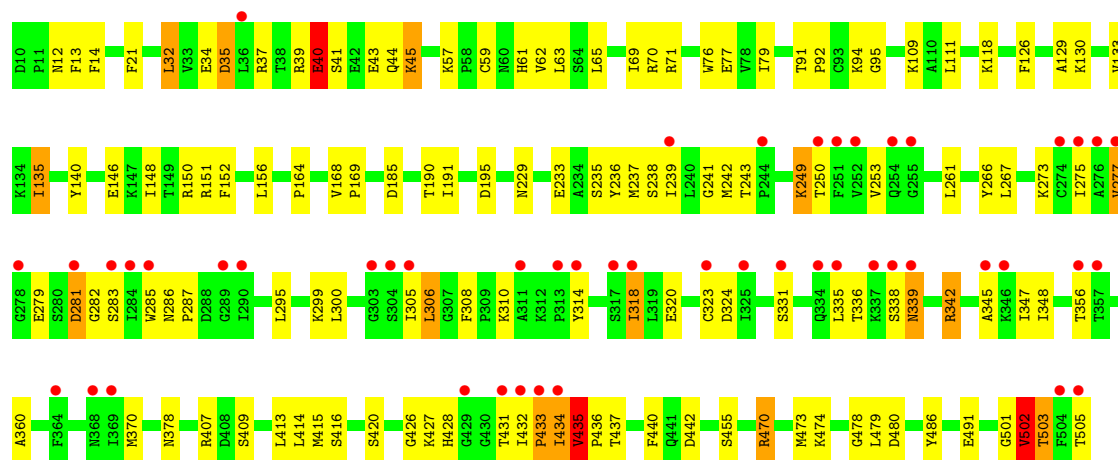


• Molecule 1: Glutamate dehydrogenase 1, mitochondrial





• Molecule 1: Glutamate dehydrogenase 1, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	96.80Å 98.38Å 124.30Å 85.94° 69.35° 61.00°	Depositor
Resolution (Å)	29.77 – 2.70 29.77 – 2.70	Depositor EDS
% Data completeness (in resolution range)	88.9 (29.77-2.70) 88.8 (29.77-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.63 (at 2.68Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.241 , 0.279 0.243 , 0.279	Depositor DCC
R_{free} test set	4609 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.455	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.008 for h,h-k,h-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	23591	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% 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: PO4

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	0.34	0/3960	0.59	3/5343 (0.1%)
1	B	0.39	0/3960	0.60	3/5343 (0.1%)
1	C	0.41	0/3960	0.67	6/5343 (0.1%)
1	D	0.39	0/3960	0.54	0/5343
1	E	0.38	0/3960	0.54	0/5343
1	F	0.37	0/3960	0.54	0/5343
All	All	0.38	0/23760	0.58	12/32058 (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	A	0	3
1	B	0	1
1	C	0	1
1	D	0	1
1	E	0	1
1	F	0	3
All	All	0	10

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	46	ARG	NE-CZ-NH2	12.18	130.16	119.20
1	C	443	ARG	NE-CZ-NH2	12.10	130.09	119.20
1	A	221	ARG	NE-CZ-NH2	11.72	129.75	119.20
1	C	46	ARG	NE-CZ-NH1	-11.70	109.80	121.50

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	ARG	NE-CZ-NH1	-11.65	109.85	121.50

There are no chirality outliers.

5 of 10 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	433	PRO	Peptide
1	A	91	THR	Mainchain,Peptide
1	B	433	PRO	Peptide
1	C	433	PRO	Peptide
1	D	433	PRO	Peptide

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	3876	0	3843	67	0
1	B	3876	0	3843	101	0
1	C	3876	0	3843	90	0
1	D	3876	0	3843	102	0
1	E	3876	0	3843	97	0
1	F	3876	0	3843	91	0
2	A	5	0	0	2	0
2	B	10	0	0	1	0
2	C	10	0	0	0	0
2	D	15	0	0	3	0
2	E	15	0	0	1	0
2	F	20	0	0	2	0
3	A	31	0	0	7	0
3	B	51	0	0	15	0
3	C	46	0	0	13	0
3	D	49	0	0	18	0
3	E	47	0	0	15	0
3	F	36	0	0	12	0
All	All	23591	0	23058	513	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 11.

The worst 5 of 513 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:370:MET:SD	3:C:746:HOH:O	2.01	1.18
1:B:227:ILE:O	3:B:702:HOH:O	1.85	0.94
1:E:58:PRO:O	3:E:701:HOH:O	1.85	0.94
1:B:315:GLU:OE2	3:B:701:HOH:O	1.84	0.93
1:F:478:GLY:N	3:F:2101:HOH:O	2.00	0.92

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	494/496 (100%)	452 (92%)	34 (7%)	8 (2%)	7	20
1	B	494/496 (100%)	453 (92%)	33 (7%)	8 (2%)	7	20
1	C	494/496 (100%)	448 (91%)	36 (7%)	10 (2%)	6	16
1	D	494/496 (100%)	452 (92%)	34 (7%)	8 (2%)	7	20
1	E	494/496 (100%)	455 (92%)	30 (6%)	9 (2%)	6	18
1	F	494/496 (100%)	448 (91%)	37 (8%)	9 (2%)	6	18
All	All	2964/2976 (100%)	2708 (91%)	204 (7%)	52 (2%)	6	18

5 of 52 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	ASP
1	A	40	GLU
1	A	434	ILE
1	B	35	ASP
1	B	40	GLU

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	413/413 (100%)	390 (94%)	23 (6%)	19	44
1	B	413/413 (100%)	390 (94%)	23 (6%)	19	44
1	C	413/413 (100%)	387 (94%)	26 (6%)	16	39
1	D	413/413 (100%)	390 (94%)	23 (6%)	19	44
1	E	413/413 (100%)	390 (94%)	23 (6%)	19	44
1	F	413/413 (100%)	389 (94%)	24 (6%)	18	42
All	All	2478/2478 (100%)	2336 (94%)	142 (6%)	18	43

5 of 142 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	470	ARG
1	F	35	ASP
1	F	318	ILE
1	C	40	GLU
1	C	35	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	232	ASN
1	D	488	ASN
1	E	302	HIS
1	E	232	ASN
1	B	232	ASN

5.3.3 RNA ⓘ

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](#)

15 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	PO4	E	601	-	4,4,4	0.92	0	6,6,6	0.39	0
2	PO4	E	603	-	4,4,4	0.80	0	6,6,6	0.51	0
2	PO4	F	2001	-	4,4,4	0.74	0	6,6,6	0.61	0
2	PO4	B	602	-	4,4,4	0.76	0	6,6,6	0.55	0
2	PO4	C	601	-	4,4,4	0.68	0	6,6,6	0.74	0
2	PO4	F	2002	-	4,4,4	0.89	0	6,6,6	0.50	0
2	PO4	F	2003	-	4,4,4	0.85	0	6,6,6	0.45	0
2	PO4	B	601	-	4,4,4	0.72	0	6,6,6	0.73	0
2	PO4	A	1001	-	4,4,4	0.86	0	6,6,6	0.64	0
2	PO4	E	602	-	4,4,4	0.96	0	6,6,6	0.44	0
2	PO4	D	601	-	4,4,4	0.89	0	6,6,6	0.33	0
2	PO4	F	2004	-	4,4,4	0.89	0	6,6,6	0.88	0
2	PO4	D	602	-	4,4,4	0.74	0	6,6,6	0.52	0
2	PO4	C	602	-	4,4,4	0.94	0	6,6,6	1.33	1 (16%)
2	PO4	D	603	-	4,4,4	1.04	0	6,6,6	0.76	0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	602	PO4	O2-P-O1	-2.35	102.63	110.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	602	PO4	1	0
2	F	2002	PO4	1	0
2	F	2003	PO4	1	0
2	A	1001	PO4	2	0
2	E	602	PO4	1	0
2	D	601	PO4	1	0
2	D	603	PO4	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	496/496 (100%)	0.51	39 (7%) 18 16	30, 65, 131, 142	0
1	B	496/496 (100%)	0.63	63 (12%) 8 7	28, 54, 103, 122	0
1	C	496/496 (100%)	1.07	113 (22%) 2 2	26, 50, 128, 153	0
1	D	496/496 (100%)	0.46	48 (9%) 13 11	25, 52, 109, 126	0
1	E	496/496 (100%)	0.15	15 (3%) 52 49	27, 49, 84, 116	0
1	F	496/496 (100%)	0.49	49 (9%) 13 10	29, 54, 113, 133	0
All	All	2976/2976 (100%)	0.55	327 (10%) 10 9	25, 54, 116, 153	0

The worst 5 of 327 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	255	GLY	11.3
1	C	285	TRP	8.4
1	C	274	CYS	8.3
1	C	290	ILE	7.6
1	C	284	ILE	7.6

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
2	PO4	B	601	5/5	0.78	0.14	75,87,101,103	0
2	PO4	C	602	5/5	0.78	0.13	39,41,54,80	0
2	PO4	F	2003	5/5	0.78	0.14	49,60,69,91	0
2	PO4	D	602	5/5	0.83	0.11	68,69,74,91	0
2	PO4	D	601	5/5	0.83	0.18	58,60,88,89	0
2	PO4	A	1001	5/5	0.85	0.12	49,65,83,100	0
2	PO4	E	603	5/5	0.86	0.16	50,62,64,78	0
2	PO4	C	601	5/5	0.91	0.08	61,65,79,81	0
2	PO4	F	2001	5/5	0.91	0.13	55,68,72,83	0
2	PO4	E	601	5/5	0.91	0.19	52,55,68,87	0
2	PO4	B	602	5/5	0.92	0.06	71,74,81,94	0
2	PO4	F	2002	5/5	0.94	0.07	55,64,70,76	0
2	PO4	E	602	5/5	0.94	0.09	53,55,63,75	0
2	PO4	F	2004	5/5	0.96	0.07	39,43,46,53	0
2	PO4	D	603	5/5	0.98	0.04	34,38,43,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.