



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 10:20 PM UTC

PDB ID : 5CDW / pdb_00005cdw
Title : Crystal Structure Analysis of a mutant Grb2 SH2 domain (W121G) with a pYVNV peptide
Authors : Papaioannou, D.; Geibel, S.; Kunze, M.; Kay, C.; Waksman, G.
Deposited on : 2015-07-05
Resolution : 2.60 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

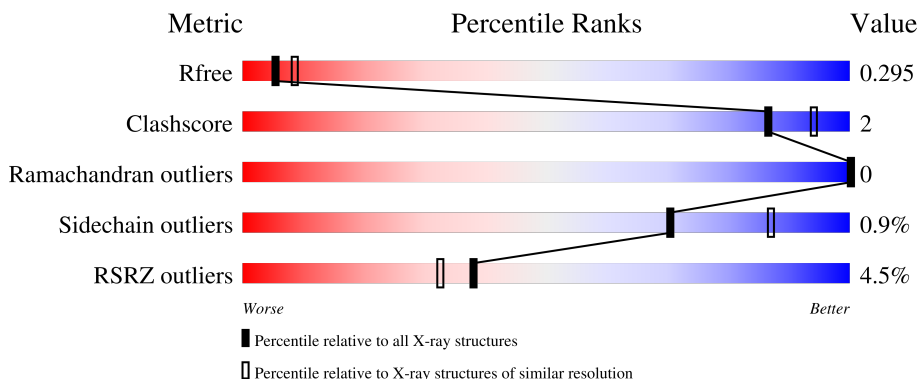
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


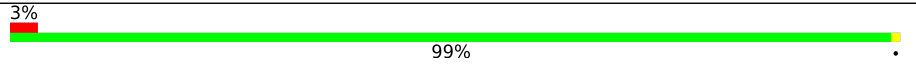
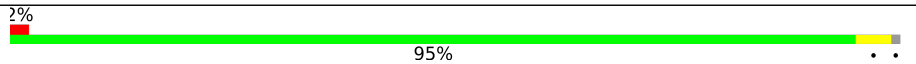
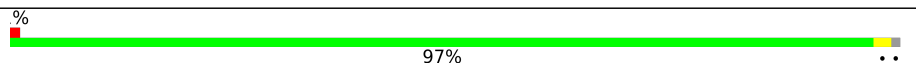
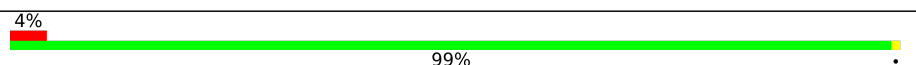
The reported resolution of this entry is 2.60 Å.

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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	100	 89% 8%
1	B	100	 99%
1	C	100	 95%
1	E	100	 97%
1	G	100	 99%



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Mol	Chain	Length	Quality of chain
1	H	100	4% 91% 6% .
1	K	100	10% 93% 7%
1	L	100	3% 93% 5% .
1	O	100	2% 95% . .
1	P	100	3% 92% . . .
1	U	100	% 95% . .
1	V	100	3% 94% . .
1	Y	100	8% 94% 5% .
1	Z	100	5% 89% 7% . .
1	c	100	7% 97% . .
1	d	100	12% 93% 5% .
2	D	6	50% 33% 17%
2	F	6	67% 17% 17%
2	I	6	17% 50% 17% 33%
2	J	6	67% 17% 17%
2	M	6	17% 50% 17% 17% 17%
2	N	6	67% 17% 17%
2	Q	6	50% 33% 17%
2	R	6	83% 17%
2	T	6	17% 67% 33%
2	X	6	50% 50%
2	a	6	83% 17%
2	b	6	67% 17% 17%
2	f	6	50% 17% 17% 17%
2	h	6	67% 17% 17%

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Mol	Chain	Length	Quality of chain
2	j	6	
2	s	6	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 25334 atoms, of which 12144 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 Growth factor receptor-bound protein 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	98	Total	C	H	N	O	S	0	0	0
			1527	500	746	140	140	1			
1	E	99	Total	C	H	N	O		0	0	0
			1494	496	720	138	140				
1	B	100	Total	C	H	N	O	S	0	0	0
			1585	516	775	145	147	2			
1	C	99	Total	C	H	N	O	S	0	0	0
			1516	500	733	139	143	1			
1	G	100	Total	C	H	N	O	S	0	0	0
			1531	504	739	142	144	2			
1	H	97	Total	C	H	N	O		0	0	0
			1475	489	714	134	138				
1	K	100	Total	C	H	N	O	S	0	0	0
			1457	490	691	134	141	1			
1	L	98	Total	C	H	N	O	S	0	0	0
			1457	489	698	132	137	1			
1	O	98	Total	C	H	N	O	S	0	0	0
			1507	496	728	140	142	1			
1	P	96	Total	C	H	N	O		0	0	0
			1466	485	711	135	135				
1	U	98	Total	C	H	N	O	S	0	0	0
			1505	496	728	138	142	1			
1	V	97	Total	C	H	N	O	S	0	0	0
			1481	490	719	136	135	1			
1	Y	99	Total	C	H	N	O	S	0	0	0
			1527	502	741	142	141	1			
1	Z	97	Total	C	H	N	O	S	0	0	0
			1457	485	702	134	135	1			
1	c	98	Total	C	H	N	O	S	0	0	0
			1413	479	669	127	137	1			
1	d	98	Total	C	H	N	O		0	0	0
			1430	482	681	130	137				

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	69	GLY	TRP	engineered mutation	UNP P62993
E	69	GLY	TRP	engineered mutation	UNP P62993
B	69	GLY	TRP	engineered mutation	UNP P62993
C	69	GLY	TRP	engineered mutation	UNP P62993
G	69	GLY	TRP	engineered mutation	UNP P62993
H	69	GLY	TRP	engineered mutation	UNP P62993
K	69	GLY	TRP	engineered mutation	UNP P62993
L	69	GLY	TRP	engineered mutation	UNP P62993
O	69	GLY	TRP	engineered mutation	UNP P62993
P	69	GLY	TRP	engineered mutation	UNP P62993
U	69	GLY	TRP	engineered mutation	UNP P62993
V	69	GLY	TRP	engineered mutation	UNP P62993
Y	69	GLY	TRP	engineered mutation	UNP P62993
Z	69	GLY	TRP	engineered mutation	UNP P62993
c	69	GLY	TRP	engineered mutation	UNP P62993
d	69	GLY	TRP	engineered mutation	UNP P62993

- Molecule 2 is a protein called SER-PTR-VAL-ASN-VAL-GLN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
2	D	5	Total	C	H	N	O	P	0	0	0
			82	26	38	6	11	1			
2	M	5	Total	C	H	N	O	P	0	0	0
			82	26	38	6	11	1			
2	s	5	Total	C	H	N	O	P	0	0	0
			79	26	35	6	11	1			
2	a	6	Total	C	H	N	O	P	0	0	0
			98	31	46	8	12	1			
2	F	5	Total	C	H	N	O	P	0	0	0
			82	26	38	6	11	1			
2	J	6	Total	C	H	N	O	P	0	0	0
			99	31	46	8	13	1			
2	N	5	Total	C	H	N	O	P	0	0	0
			82	26	38	6	11	1			
2	R	6	Total	C	H	N	O	P	0	0	0
			99	31	46	8	13	1			
2	T	6	Total	C	H	N	O	P	0	0	0
			89	29	40	7	12	1			
2	X	6	Total	C	H	N	O	P	0	0	0
			89	29	40	7	12	1			
2	b	5	Total	C	H	N	O	P	0	0	0
			82	26	38	6	11	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				P
2	f	5	Total 82	C 26	H 38	N 6	O 11	P 1	0	0	0
2	h	6	Total 99	C 31	H 46	N 8	O 13	P 1	0	0	0
2	j	5	Total 82	C 26	H 38	N 6	O 11	P 1	0	0	0
2	I	6	Total 99	C 31	H 46	N 8	O 13	P 1	0	0	0
2	Q	5	Total 82	C 26	H 38	N 6	O 11	P 1	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total 4	O 4	0	0
3	E	9	Total 9	O 9	0	0
3	B	5	Total 5	O 5	0	0
3	C	10	Total 10	O 10	0	0
3	G	3	Total 3	O 3	0	0
3	H	4	Total 4	O 4	0	0
3	K	3	Total 3	O 3	0	0
3	L	5	Total 5	O 5	0	0
3	O	7	Total 7	O 7	0	0
3	P	12	Total 12	O 12	0	0
3	U	6	Total 6	O 6	0	0
3	V	6	Total 6	O 6	0	0
3	Y	7	Total 7	O 7	0	0
3	Z	7	Total 7	O 7	0	0

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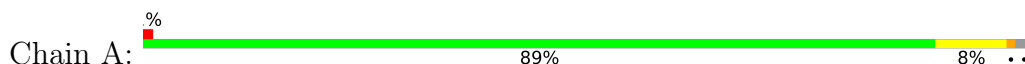
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	c	4	Total O 4 4	0	0
3	d	4	Total O 4 4	0	0
3	b	1	Total O 1 1	0	0
3	f	1	Total O 1 1	0	0
3	j	1	Total O 1 1	0	0

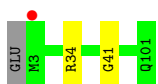
3 Residue-property plots [i](#)

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.

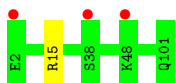
- Molecule 1: Growth factor receptor-bound protein 2



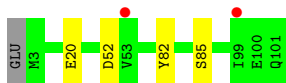
- Molecule 1: Growth factor receptor-bound protein 2



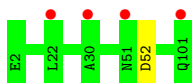
- Molecule 1: Growth factor receptor-bound protein 2



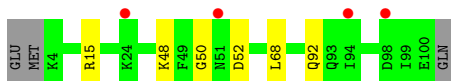
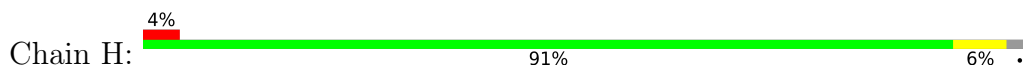
- Molecule 1: Growth factor receptor-bound protein 2



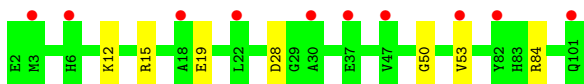
- Molecule 1: Growth factor receptor-bound protein 2



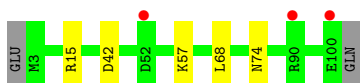
- Molecule 1: Growth factor receptor-bound protein 2



- Molecule 1: Growth factor receptor-bound protein 2



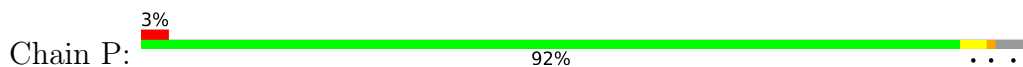
- Molecule 1: Growth factor receptor-bound protein 2



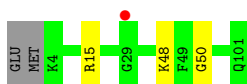
- Molecule 1: Growth factor receptor-bound protein 2



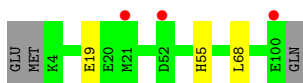
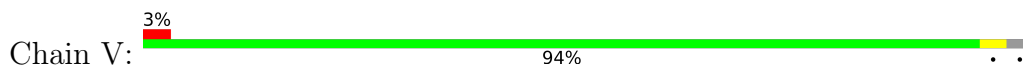
- Molecule 1: Growth factor receptor-bound protein 2



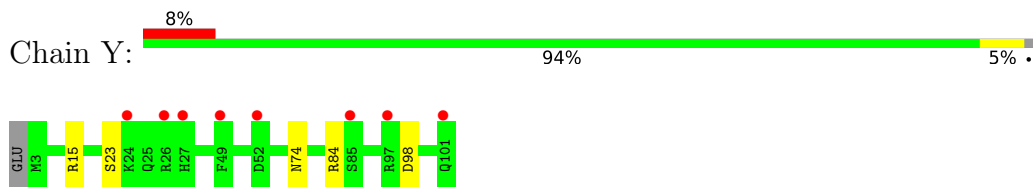
- Molecule 1: Growth factor receptor-bound protein 2



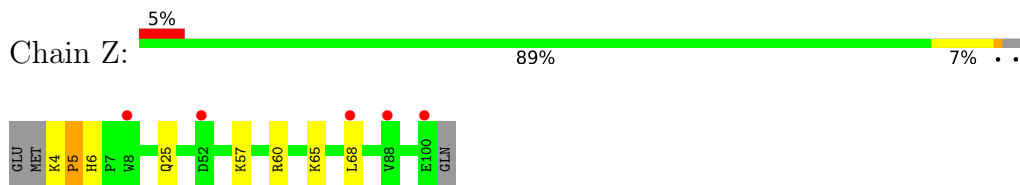
- Molecule 1: Growth factor receptor-bound protein 2



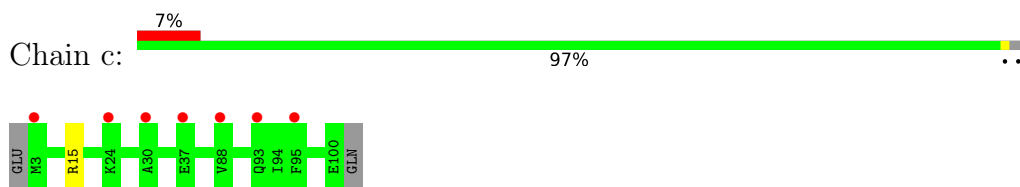
- Molecule 1: Growth factor receptor-bound protein 2



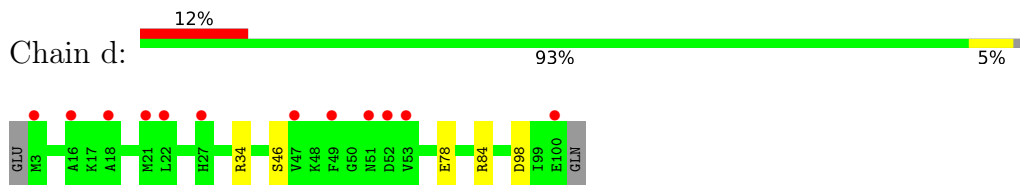
- Molecule 1: Growth factor receptor-bound protein 2



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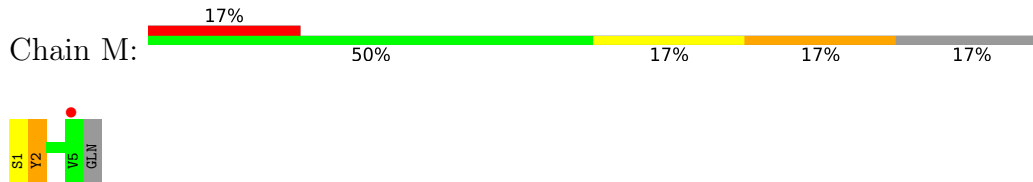
- Molecule 1: Growth factor receptor-bound protein 2



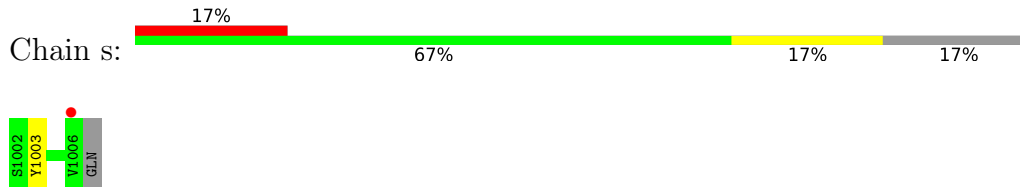
- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN




- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN

Chain a:  83% 17%



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN

Chain F:  67% 17% 17%



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN

Chain J:  67% 17% 17%




- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN

Chain N:  67% 17% 17%



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN

Chain R:  83% 17%



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN

Chain T:  17% 67% 33%



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN

Chain X:  50% 50%



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN

Chain b:  67% 17% 17%



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN



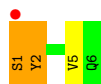
- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN



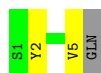
- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN



- Molecule 2: SER-PTR-VAL-ASN-VAL-GLN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.65Å 72.14Å 172.14Å 90.00° 93.31° 90.00°	Depositor
Resolution (Å)	41.26 – 2.60 41.26 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.0 (41.26-2.60) 99.2 (41.26-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.61Å)	Xtrriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.236 , 0.294 0.246 , 0.295	Depositor DCC
R_{free} test set	3103 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	46.0	Xtrriage
Anisotropy	1.023	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	25334	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.64 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.5401e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.27	0/801	0.62	0/1078
1	B	0.27	0/830	0.58	0/1115
1	C	0.26	0/803	0.57	0/1083
1	E	0.26	0/794	0.56	0/1072
1	G	0.26	0/812	0.59	0/1095
1	H	0.25	0/781	0.57	0/1054
1	K	0.27	0/786	0.56	0/1065
1	L	0.27	0/779	0.57	0/1053
1	O	0.26	0/799	0.58	0/1078
1	P	0.27	0/775	0.58	0/1045
1	U	0.26	0/797	0.56	0/1075
1	V	0.27	0/782	0.59	0/1054
1	Y	0.25	0/806	0.59	0/1086
1	Z	0.28	0/775	0.66	1/1047 (0.1%)
1	c	0.27	0/764	0.57	0/1036
1	d	0.27	0/769	0.59	0/1042
2	D	0.13	0/26	0.51	0/33
2	F	0.23	0/26	0.62	0/33
2	I	0.20	0/35	0.43	0/45
2	J	0.24	0/35	0.47	0/45
2	M	0.25	0/26	0.72	0/33
2	N	0.19	0/26	0.43	0/33
2	Q	0.20	0/26	0.49	0/33
2	R	0.16	0/35	0.40	0/45
2	T	0.17	0/31	0.40	0/40
2	X	0.22	0/31	0.43	0/40
2	a	0.19	0/34	0.48	0/44
2	b	0.16	0/26	0.42	0/33
2	f	0.18	0/26	0.41	0/33
2	h	0.22	0/35	0.47	0/45
2	j	0.21	0/26	0.52	0/33
2	s	0.23	0/26	0.57	0/33

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
All	All	0.26	0/13123	0.58	1/17679 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Z	5	PRO	CA-C-O	-5.13	114.92	122.31

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	781	746	750	8	0
1	B	810	775	779	1	0
1	C	783	733	737	3	0
1	E	774	720	724	2	0
1	G	792	739	743	1	0
1	H	761	714	718	3	0
1	K	766	691	695	5	0
1	L	759	698	702	3	0
1	O	779	728	732	2	0
1	P	755	711	715	4	0
1	U	777	728	732	2	0
1	V	762	719	723	1	0
1	Y	786	741	745	4	0
1	Z	755	702	706	6	0
1	c	744	669	673	1	0
1	d	749	681	685	4	0
2	D	44	38	38	1	0
2	F	44	38	38	1	0
2	I	53	46	46	4	0
2	J	53	46	46	1	0
2	M	44	38	38	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	N	44	38	39	0	0
2	Q	44	38	38	0	0
2	R	53	46	46	0	0
2	T	49	40	40	1	0
2	X	49	40	41	2	0
2	a	52	46	46	1	0
2	b	44	38	38	0	0
2	f	44	38	38	2	0
2	h	53	46	46	1	0
2	j	44	38	38	1	0
2	s	44	35	35	0	0
3	A	4	0	0	1	0
3	B	5	0	0	0	0
3	C	10	0	0	2	0
3	E	9	0	0	1	0
3	G	3	0	0	1	0
3	H	4	0	0	0	0
3	K	3	0	0	1	0
3	L	5	0	0	1	0
3	O	7	0	0	0	0
3	P	12	0	0	3	0
3	U	6	0	0	1	0
3	V	6	0	0	0	0
3	Y	7	0	0	1	0
3	Z	7	0	0	4	0
3	b	1	0	0	0	0
3	c	4	0	0	0	0
3	d	4	0	0	0	0
3	f	1	0	0	1	0
3	j	1	0	0	0	0
All	All	13190	12144	12210	53	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 2.

The worst 5 of 53 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:Z:4:LYS:N	3:Z:201:HOH:O	1.96	0.97
1:A:71:VAL:O	3:A:201:HOH:O	1.87	0.93
2:f:1:SER:OG	3:f:101:HOH:O	1.95	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:68:LEU:O	3:P:201:HOH:O	1.96	0.81
1:Z:4:LYS:O	3:Z:201:HOH:O	1.97	0.81

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	96/100 (96%)	94 (98%)	2 (2%)	0	100	100
1	B	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
1	C	97/100 (97%)	94 (97%)	3 (3%)	0	100	100
1	E	97/100 (97%)	96 (99%)	1 (1%)	0	100	100
1	G	98/100 (98%)	92 (94%)	6 (6%)	0	100	100
1	H	95/100 (95%)	92 (97%)	3 (3%)	0	100	100
1	K	98/100 (98%)	91 (93%)	7 (7%)	0	100	100
1	L	96/100 (96%)	92 (96%)	4 (4%)	0	100	100
1	O	96/100 (96%)	93 (97%)	3 (3%)	0	100	100
1	P	94/100 (94%)	92 (98%)	2 (2%)	0	100	100
1	U	96/100 (96%)	96 (100%)	0	0	100	100
1	V	95/100 (95%)	95 (100%)	0	0	100	100
1	Y	97/100 (97%)	94 (97%)	3 (3%)	0	100	100
1	Z	95/100 (95%)	91 (96%)	4 (4%)	0	100	100
1	c	96/100 (96%)	92 (96%)	4 (4%)	0	100	100
1	d	96/100 (96%)	91 (95%)	5 (5%)	0	100	100
2	D	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	F	2/6 (33%)	1 (50%)	1 (50%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	I	3/6 (50%)	3 (100%)	0	0	100	100
2	J	3/6 (50%)	3 (100%)	0	0	100	100
2	M	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	N	2/6 (33%)	2 (100%)	0	0	100	100
2	Q	2/6 (33%)	2 (100%)	0	0	100	100
2	R	3/6 (50%)	3 (100%)	0	0	100	100
2	T	3/6 (50%)	3 (100%)	0	0	100	100
2	X	3/6 (50%)	3 (100%)	0	0	100	100
2	a	3/6 (50%)	2 (67%)	1 (33%)	0	100	100
2	b	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	f	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	h	3/6 (50%)	2 (67%)	1 (33%)	0	100	100
2	j	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	s	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
All	All	1579/1696 (93%)	1521 (96%)	58 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	81/88 (92%)	80 (99%)	1 (1%)	63	83
1	B	85/88 (97%)	85 (100%)	0	100	100
1	C	81/88 (92%)	80 (99%)	1 (1%)	63	83
1	E	78/88 (89%)	78 (100%)	0	100	100
1	G	81/88 (92%)	81 (100%)	0	100	100
1	H	78/88 (89%)	76 (97%)	2 (3%)	40	68
1	K	75/88 (85%)	75 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	75/88 (85%)	74 (99%)	1 (1%)	61	82
1	O	80/88 (91%)	80 (100%)	0	100	100
1	P	77/88 (88%)	76 (99%)	1 (1%)	61	82
1	U	80/88 (91%)	80 (100%)	0	100	100
1	V	78/88 (89%)	77 (99%)	1 (1%)	61	82
1	Y	80/88 (91%)	80 (100%)	0	100	100
1	Z	76/88 (86%)	74 (97%)	2 (3%)	40	68
1	c	72/88 (82%)	72 (100%)	0	100	100
1	d	74/88 (84%)	74 (100%)	0	100	100
2	D	4/5 (80%)	4 (100%)	0	100	100
2	F	4/5 (80%)	4 (100%)	0	100	100
2	I	5/5 (100%)	4 (80%)	1 (20%)	1	2
2	J	5/5 (100%)	5 (100%)	0	100	100
2	M	4/5 (80%)	4 (100%)	0	100	100
2	N	4/5 (80%)	4 (100%)	0	100	100
2	Q	4/5 (80%)	3 (75%)	1 (25%)	0	1
2	R	5/5 (100%)	5 (100%)	0	100	100
2	T	4/5 (80%)	4 (100%)	0	100	100
2	X	4/5 (80%)	4 (100%)	0	100	100
2	a	5/5 (100%)	5 (100%)	0	100	100
2	b	4/5 (80%)	4 (100%)	0	100	100
2	f	4/5 (80%)	4 (100%)	0	100	100
2	h	5/5 (100%)	5 (100%)	0	100	100
2	j	4/5 (80%)	3 (75%)	1 (25%)	0	1
2	s	4/5 (80%)	4 (100%)	0	100	100
All	All	1320/1488 (89%)	1308 (99%)	12 (1%)	70	87

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	Z	25	GLN
1	Z	68	LEU
2	Q	5	VAL

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Mol	Chain	Res	Type
2	j	1	SER
1	H	92	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	H	91	ASN
1	c	91	ASN
1	K	54	GLN
1	c	92	GLN
1	Y	6	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](#)

16 non-standard protein/DNA/RNA residues 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	PTR	X	2	2	15,16,17	1.23	1 (6%)	17,22,24	0.57	0
2	PTR	j	2	2	15,16,17	1.36	1 (6%)	17,22,24	0.48	0
2	PTR	s	1003	2	15,16,17	1.35	1 (6%)	17,22,24	0.54	0
2	PTR	R	2	2	15,16,17	1.34	1 (6%)	17,22,24	0.52	0
2	PTR	T	2	2	15,16,17	1.22	1 (6%)	17,22,24	0.56	0
2	PTR	f	2	2	15,16,17	1.23	1 (6%)	17,22,24	0.60	0
2	PTR	N	2	2	15,16,17	1.22	1 (6%)	17,22,24	0.55	0
2	PTR	I	2	2	15,16,17	1.25	1 (6%)	17,22,24	0.60	0
2	PTR	D	2	2	15,16,17	1.35	1 (6%)	17,22,24	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PTR	b	2	2	15,16,17	1.34	1 (6%)	17,22,24	0.52	0
2	PTR	M	2	2	15,16,17	1.34	1 (6%)	17,22,24	0.53	0
2	PTR	J	2	2	15,16,17	1.33	1 (6%)	17,22,24	0.53	0
2	PTR	a	2	2	15,16,17	1.33	1 (6%)	17,22,24	0.50	0
2	PTR	F	2	2	15,16,17	1.37	1 (6%)	17,22,24	0.57	0
2	PTR	h	2	2	15,16,17	1.37	1 (6%)	17,22,24	0.55	0
2	PTR	Q	2	2	15,16,17	1.26	1 (6%)	17,22,24	0.59	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	PTR	X	2	2	-	1/10/11/13	0/1/1/1
2	PTR	j	2	2	-	1/10/11/13	0/1/1/1
2	PTR	s	1003	2	-	0/10/11/13	0/1/1/1
2	PTR	R	2	2	-	1/10/11/13	0/1/1/1
2	PTR	T	2	2	-	1/10/11/13	0/1/1/1
2	PTR	f	2	2	-	1/10/11/13	0/1/1/1
2	PTR	N	2	2	-	2/10/11/13	0/1/1/1
2	PTR	I	2	2	-	0/10/11/13	0/1/1/1
2	PTR	D	2	2	-	0/10/11/13	0/1/1/1
2	PTR	b	2	2	-	0/10/11/13	0/1/1/1
2	PTR	M	2	2	-	0/10/11/13	0/1/1/1
2	PTR	J	2	2	-	1/10/11/13	0/1/1/1
2	PTR	a	2	2	-	0/10/11/13	0/1/1/1
2	PTR	F	2	2	-	0/10/11/13	0/1/1/1
2	PTR	h	2	2	-	1/10/11/13	0/1/1/1
2	PTR	Q	2	2	-	1/10/11/13	0/1/1/1

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	PTR	OH-CZ	-4.64	1.30	1.40
2	h	2	PTR	OH-CZ	-4.63	1.30	1.40
2	Q	2	PTR	OH-CZ	-4.60	1.30	1.40
2	I	2	PTR	OH-CZ	-4.57	1.30	1.40
2	j	2	PTR	OH-CZ	-4.54	1.30	1.40

There are no bond angle outliers.

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	N	2	PTR	O-C-CA-CB
2	R	2	PTR	O-C-CA-CB
2	h	2	PTR	O-C-CA-CB
2	j	2	PTR	O-C-CA-CB
2	T	2	PTR	CZ-OH-P-O1P

There are no ring outliers.

7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	f	2	PTR	1	0
2	I	2	PTR	3	0
2	M	2	PTR	2	0
2	J	2	PTR	1	0
2	a	2	PTR	1	0
2	F	2	PTR	1	0
2	h	2	PTR	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	98/100 (98%)	0.51	1 (1%) 79 76	49, 58, 68, 72	0
1	B	100/100 (100%)	0.56	3 (3%) 52 47	49, 57, 73, 122	0
1	C	99/100 (99%)	0.60	2 (2%) 65 60	49, 60, 74, 78	0
1	E	99/100 (99%)	0.52	1 (1%) 79 76	48, 59, 72, 84	0
1	G	100/100 (100%)	0.77	4 (4%) 42 37	50, 69, 85, 96	0
1	H	97/100 (97%)	0.66	4 (4%) 41 36	51, 65, 80, 86	0
1	K	100/100 (100%)	1.04	10 (10%) 12 9	57, 68, 88, 98	0
1	L	98/100 (98%)	0.69	3 (3%) 51 45	56, 66, 77, 88	0
1	O	98/100 (98%)	0.47	2 (2%) 65 60	48, 57, 66, 79	0
1	P	96/100 (96%)	0.56	3 (3%) 51 45	48, 57, 66, 70	0
1	U	98/100 (98%)	0.64	1 (1%) 79 76	53, 66, 78, 86	0
1	V	97/100 (97%)	0.74	3 (3%) 51 45	52, 68, 78, 82	0
1	Y	99/100 (99%)	0.72	8 (8%) 18 14	52, 60, 73, 78	0
1	Z	97/100 (97%)	0.83	5 (5%) 33 27	53, 64, 77, 83	0
1	c	98/100 (98%)	0.85	7 (7%) 22 17	58, 69, 81, 87	0
1	d	98/100 (98%)	0.85	12 (12%) 8 6	55, 65, 80, 85	0
2	D	4/6 (66%)	1.23	0 100 100	57, 58, 61, 62	0
2	F	4/6 (66%)	1.44	0 100 100	61, 63, 65, 70	0
2	I	5/6 (83%)	1.26	1 (20%) 3 2	64, 66, 76, 86	0
2	J	5/6 (83%)	0.74	0 100 100	59, 61, 69, 70	0
2	M	4/6 (66%)	0.83	1 (25%) 2 1	72, 75, 79, 84	0
2	N	4/6 (66%)	1.46	0 100 100	60, 63, 64, 77	0
2	Q	4/6 (66%)	1.15	0 100 100	67, 72, 73, 77	0
2	R	5/6 (83%)	0.55	0 100 100	58, 62, 68, 70	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
2	T	5/6 (83%)	1.06	1 (20%) 3 2	57, 63, 66, 71	0
2	X	5/6 (83%)	0.90	0 100 100	60, 63, 68, 70	0
2	a	5/6 (83%)	0.95	0 100 100	56, 58, 60, 64	0
2	b	4/6 (66%)	0.45	0 100 100	55, 58, 63, 66	0
2	f	4/6 (66%)	0.57	0 100 100	53, 58, 60, 63	0
2	h	5/6 (83%)	0.97	0 100 100	70, 74, 83, 95	0
2	j	4/6 (66%)	1.54	1 (25%) 2 1	68, 73, 74, 75	0
2	s	4/6 (66%)	1.22	1 (25%) 2 1	57, 59, 65, 66	0
All	All	1643/1696 (96%)	0.70	74 (4%) 38 32	48, 63, 80, 122	0

The worst 5 of 74 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	101	GLN	3.9
1	L	90	ARG	3.9
1	c	30	ALA	3.9
1	d	52	ASP	3.9
1	L	100	GLU	3.7

6.2 Non-standard residues in protein, DNA, RNA chains [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	PTR	F	2	16/17	0.88	0.13	58,60,72,72	0
2	PTR	Q	2	16/17	0.88	0.12	64,66,78,81	0
2	PTR	h	2	16/17	0.89	0.13	61,66,79,81	0
2	PTR	J	2	16/17	0.90	0.12	56,59,70,71	0
2	PTR	j	2	16/17	0.91	0.12	60,63,75,76	0
2	PTR	M	2	16/17	0.91	0.11	62,70,84,86	0
2	PTR	I	2	16/17	0.92	0.12	53,58,70,71	0
2	PTR	b	2	16/17	0.92	0.11	54,57,69,69	0
2	PTR	a	2	16/17	0.93	0.11	50,52,62,64	0
2	PTR	N	2	16/17	0.93	0.10	58,63,74,76	0
2	PTR	X	2	16/17	0.93	0.11	55,57,69,69	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PTR	D	2	16/17	0.93	0.11	47,51,61,63	0
2	PTR	R	2	16/17	0.94	0.08	53,57,68,69	0
2	PTR	f	2	16/17	0.94	0.10	51,55,65,65	0
2	PTR	T	2	16/17	0.95	0.11	52,55,67,68	0
2	PTR	s	1003	16/17	0.96	0.09	50,53,64,64	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.