



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:28 AM UTC

PDB ID : 3UYC / pdb_00003uyc
Title : Designed protein KE59 R8_2/7A
Authors : Khersonsky, O.; Kiss, G.; Roethlisberger, D.; Dym, O.; Albeck, S.; Houk, K.N.; Baker, D.; Tawfik, D.S.; Israel Structural Proteomics Center (ISPC)
Deposited on : 2011-12-06
Resolution : 2.20 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

<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

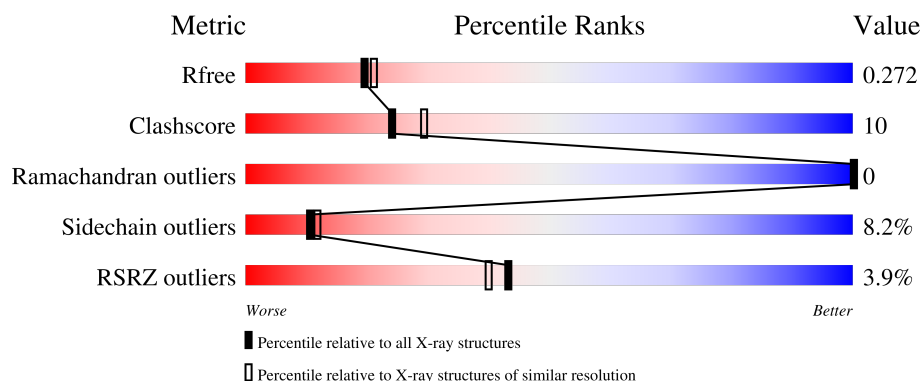
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

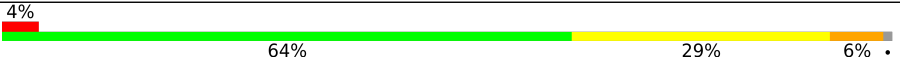
The reported resolution of this entry is 2.20 Å.

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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	249	
1	B	249	

2 Entry composition [i](#)

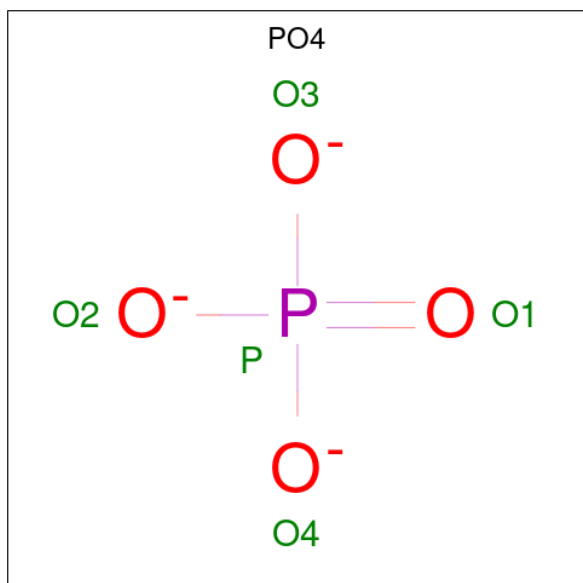
There are 3 unique types of molecules in this entry. The entry contains 3904 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Kemp eliminase KE59 R8_2/7A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	0	0	0
			1883	1206	323	349	5			
1	B	246	Total	C	N	O	S	0	0	0
			1954	1245	340	364	5			

- 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		

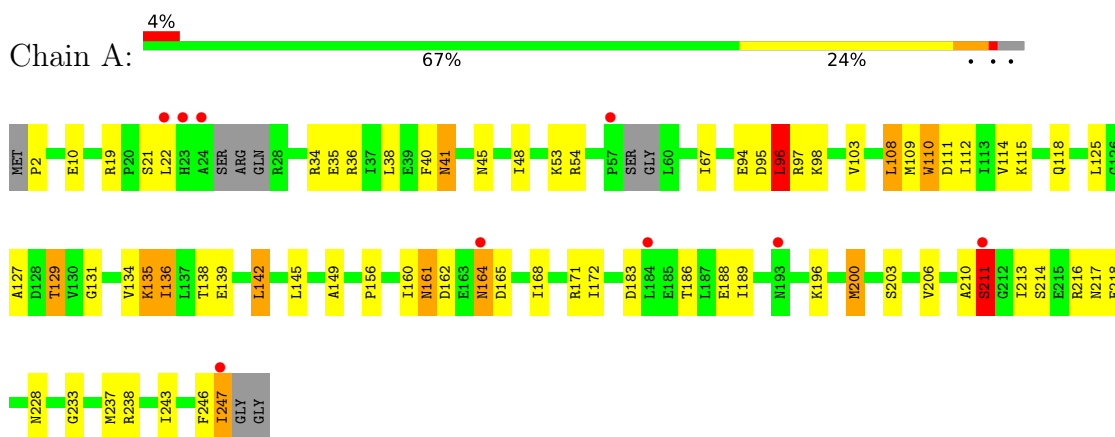
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	23	Total 23	O 23	0	0
3	B	34	Total 34	O 34	0	0

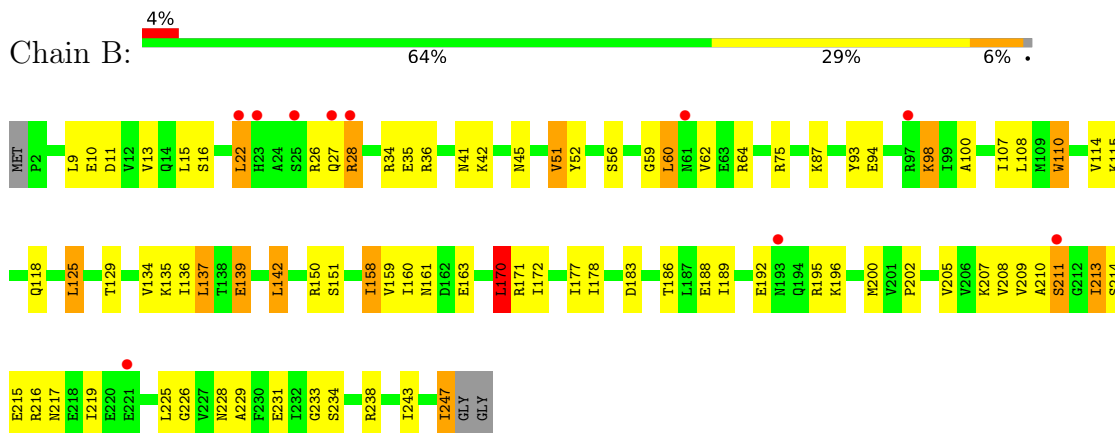
3 Residue-property plots

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: Kemp eliminase KE59 R8_2/7A



• Molecule 1: Kemp eliminase KE59 R8_2/7A



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	99.57Å 65.47Å 75.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.59 – 2.20 41.59 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.9 (41.59-2.20) 97.9 (41.59-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.66 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.4.0067	Depositor
R, R_{free}	0.209 , 0.275 0.217 , 0.272	Depositor DCC
R_{free} test set	1282 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 26.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3904	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 66.25 % 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 6.2248e-06. 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

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	1.65	18/1908 (0.9%)	1.40	15/2576 (0.6%)
1	B	1.65	19/1982 (1.0%)	1.44	18/2673 (0.7%)
All	All	1.65	37/3890 (1.0%)	1.42	33/5249 (0.6%)

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	2
1	B	0	1
All	All	0	3

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	136	ILE	CA-CB	8.05	1.65	1.54
1	A	48	ILE	CA-CB	7.25	1.62	1.54
1	A	188	GLU	C-O	-7.23	1.15	1.23
1	A	41	ASN	C-O	-7.17	1.15	1.24
1	B	36	ARG	C-O	-6.74	1.16	1.24
1	B	129	THR	C-O	6.66	1.31	1.23
1	B	107	ILE	C-O	-6.31	1.17	1.24
1	A	217	ASN	C-O	-6.23	1.16	1.24
1	A	36	ARG	C-O	-6.16	1.17	1.24
1	A	35	GLU	C-O	-6.09	1.17	1.24
1	B	158	ILE	CA-C	6.07	1.60	1.52
1	B	216	ARG	C-O	-6.07	1.17	1.24
1	B	209	VAL	C-O	-6.01	1.17	1.24
1	B	35	GLU	C-O	-6.00	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	129	THR	C-O	-5.99	1.16	1.23
1	B	52	TYR	C-O	-5.92	1.16	1.24
1	A	109	MET	C-O	-5.90	1.17	1.24
1	A	189	ILE	C-O	-5.83	1.17	1.24
1	A	34	ARG	C-O	-5.79	1.17	1.24
1	B	189	ILE	C-O	-5.66	1.18	1.24
1	B	219	ILE	CA-CB	5.61	1.60	1.54
1	B	178	ILE	CA-CB	5.56	1.60	1.54
1	B	139	GLU	C-O	-5.54	1.17	1.24
1	A	200	MET	N-CA	5.54	1.53	1.46
1	B	56	SER	C-O	-5.44	1.17	1.24
1	A	40	PHE	C-O	-5.41	1.17	1.24
1	B	163	GLU	N-CA	5.34	1.52	1.46
1	A	135	LYS	C-O	-5.34	1.17	1.24
1	A	206	VAL	N-CA	5.34	1.52	1.46
1	A	38	LEU	C-O	-5.21	1.18	1.24
1	B	34	ARG	C-O	-5.19	1.18	1.24
1	B	188	GLU	C-O	-5.17	1.17	1.23
1	A	218	GLU	C-O	-5.14	1.18	1.24
1	B	75	ARG	C-O	-5.11	1.17	1.24
1	B	170	LEU	C-O	-5.09	1.18	1.24
1	B	238	ARG	C-O	-5.04	1.18	1.24
1	A	168	ILE	CA-CB	5.00	1.60	1.54

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	59	GLY	N-CA-C	11.25	126.95	113.79
1	A	164	ASN	N-CA-C	8.09	119.72	111.07
1	B	211	SER	N-CA-C	7.69	120.96	108.34
1	B	161	ASN	N-CA-CB	-7.32	101.26	111.00
1	B	163	GLU	N-CA-C	7.31	119.24	111.28
1	B	108	LEU	CD1-CG-CD2	-6.92	95.57	110.80
1	B	26	ARG	CB-CA-C	-6.84	99.24	109.90
1	B	108	LEU	CA-CB-CG	6.66	139.61	116.30
1	B	214	SER	CB-CA-C	-6.22	101.14	109.16
1	A	19	ARG	CA-C-N	-5.94	113.64	119.76
1	A	19	ARG	C-N-CA	-5.94	113.64	119.76
1	A	35	GLU	N-CA-C	5.90	118.19	111.11
1	A	96	LEU	CD1-CG-CD2	-5.85	97.94	110.80
1	A	145	LEU	CD1-CG-CD2	-5.72	98.22	110.80
1	A	203	SER	N-CA-C	5.55	118.09	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	213	ILE	N-CA-C	5.54	116.20	108.89
1	A	211	SER	N-CA-C	-5.52	102.02	110.14
1	B	27	GLN	CB-CA-C	-5.51	101.52	110.79
1	B	135	LYS	N-CA-C	5.41	119.36	112.87
1	B	51	VAL	N-CA-C	5.35	116.29	108.58
1	B	234	SER	N-CA-C	5.35	116.80	111.07
1	A	34	ARG	N-CA-C	5.30	117.47	111.11
1	A	103	VAL	CA-C-N	-5.23	114.28	122.49
1	A	103	VAL	C-N-CA	-5.23	114.28	122.49
1	A	135	LYS	N-CA-C	5.22	119.14	112.87
1	B	26	ARG	NE-CZ-NH1	-5.19	116.31	121.50
1	B	211	SER	CA-C-N	-5.18	111.25	121.41
1	B	211	SER	C-N-CA	-5.18	111.25	121.41
1	B	15	LEU	CB-CA-C	-5.10	102.32	110.79
1	A	233	GLY	CA-C-N	5.03	128.36	120.82
1	A	233	GLY	C-N-CA	5.03	128.36	120.82
1	B	217	ASN	N-CA-C	5.02	117.40	111.33
1	A	138	THR	N-CA-C	-5.01	104.07	110.53

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	ALA	Peptide
1	A	211	SER	Peptide
1	B	210	ALA	Peptide

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	1883	0	1922	38	0
1	B	1954	0	2016	41	0
2	A	5	0	0	0	0
2	B	5	0	0	1	0
3	A	23	0	0	0	0
3	B	34	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3904	0	3938	78	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 10.

All (78) 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:A:211:SER:HB2	1:A:213:ILE:HG13	1.25	1.11
1:A:211:SER:CB	1:A:213:ILE:HG13	1.88	1.04
1:A:211:SER:HB2	1:A:213:ILE:CG1	1.99	0.92
1:B:60:LEU:CD2	1:B:62:VAL:HG23	2.01	0.90
1:B:211:SER:O	2:B:301:PO4:O3	1.97	0.82
1:A:114:VAL:H	1:A:118:GLN:HE22	1.34	0.74
1:B:114:VAL:H	1:B:118:GLN:HE22	1.34	0.73
1:B:136:ILE:HG22	1:B:137:LEU:HD13	1.74	0.69
1:B:114:VAL:H	1:B:118:GLN:NE2	1.92	0.67
1:A:211:SER:HB2	1:A:213:ILE:H	1.62	0.65
1:A:54:ARG:NH2	1:A:95:ASP:OD2	2.30	0.65
1:B:243:ILE:HD11	1:B:247:ILE:HD11	1.81	0.62
1:A:211:SER:HB2	1:A:213:ILE:CB	2.30	0.62
1:B:60:LEU:HD21	1:B:62:VAL:HG23	1.81	0.62
1:A:211:SER:HB2	1:A:213:ILE:N	2.14	0.62
1:A:108:LEU:HD13	1:A:129:THR:HG23	1.82	0.61
1:A:183:ASP:OD2	1:A:186:THR:HG23	2.00	0.60
1:B:60:LEU:HD22	1:B:62:VAL:HG23	1.79	0.60
1:A:114:VAL:H	1:A:118:GLN:NE2	1.97	0.60
1:B:170:LEU:HD21	1:B:202:PRO:HD3	1.84	0.59
1:B:211:SER:HB2	1:B:231:GLU:O	2.04	0.58
1:A:211:SER:CB	1:A:213:ILE:CG1	2.70	0.57
1:A:139:GLU:OE2	1:A:171:ARG:HD2	2.02	0.57
1:B:139:GLU:OE2	1:B:171:ARG:HD2	2.05	0.56
1:B:22:LEU:HD22	1:B:125:LEU:HD11	1.87	0.56
1:A:164:ASN:O	1:A:164:ASN:OD1	2.23	0.55
1:A:142:LEU:HD13	1:A:172:ILE:CD1	2.37	0.55
1:A:67:ILE:HD11	1:A:98:LYS:HB3	1.89	0.54
1:A:243:ILE:HD11	1:A:247:ILE:HD11	1.90	0.53
1:B:142:LEU:HD13	1:B:172:ILE:CD1	2.39	0.53
1:B:60:LEU:HD21	1:B:62:VAL:CG2	2.38	0.53
1:A:164:ASN:OD1	1:A:164:ASN:C	2.52	0.53
1:B:41:ASN:ND2	1:B:228:ASN:HD22	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:ARG:NH1	1:A:246:PHE:O	2.42	0.51
1:A:196:LYS:O	1:A:200:MET:HG3	2.10	0.51
1:A:135:LYS:NZ	1:A:162:ASP:OD2	2.43	0.51
1:A:111:ASP:OD1	1:A:112:ILE:N	2.42	0.51
1:B:243:ILE:O	1:B:247:ILE:HD13	2.10	0.51
1:B:183:ASP:OD2	1:B:186:THR:HG23	2.12	0.50
1:A:211:SER:CB	1:A:213:ILE:H	2.25	0.49
1:A:53:LYS:O	1:A:237:MET:HE2	2.12	0.49
1:A:142:LEU:HD13	1:A:172:ILE:HD11	1.94	0.49
1:B:150:ARG:HD2	3:B:268:HOH:O	2.12	0.49
1:B:211:SER:CB	1:B:231:GLU:O	2.61	0.49
1:B:134:VAL:HG22	1:B:142:LEU:HD11	1.94	0.48
1:B:94:GLU:O	1:B:98:LYS:HG2	2.12	0.48
1:A:96:LEU:HD11	1:A:127:ALA:HB2	1.96	0.48
1:A:183:ASP:OD2	1:A:186:THR:CG2	2.62	0.47
1:B:196:LYS:O	1:B:200:MET:HG3	2.15	0.47
1:A:149:ALA:HB3	1:A:156:PRO:HG3	1.96	0.47
1:B:110:TRP:HA	1:B:110:TRP:CE3	2.50	0.47
1:A:41:ASN:ND2	1:A:228:ASN:HD22	2.13	0.46
1:A:134:VAL:HG11	1:A:165:ASP:HA	1.97	0.46
1:B:207:LYS:HE2	1:B:226:GLY:O	2.15	0.46
1:B:208:VAL:HG22	1:B:229:ALA:HB3	1.97	0.46
1:B:243:ILE:CD1	1:B:247:ILE:HD11	2.46	0.46
1:B:16:SER:O	1:B:115:LYS:HE2	2.15	0.46
1:A:238:ARG:NH1	1:B:151:SER:OG	2.47	0.45
1:B:11:ASP:OD2	1:B:87:LYS:NZ	2.48	0.43
1:A:110:TRP:CE3	1:A:110:TRP:HA	2.53	0.43
1:B:28:ARG:NH1	1:B:100:ALA:O	2.51	0.43
1:B:9:LEU:HD23	1:B:9:LEU:HA	1.89	0.43
1:B:51:VAL:HB	1:B:233:GLY:HA2	2.01	0.42
1:A:115:LYS:H	1:A:118:GLN:HE21	1.67	0.42
1:B:177:ILE:HD12	1:B:205:VAL:HG11	2.00	0.42
1:A:108:LEU:HD11	1:A:131:GLY:N	2.35	0.41
1:B:22:LEU:HD22	1:B:125:LEU:CD1	2.50	0.41
1:B:13:VAL:HG13	1:B:114:VAL:HG11	2.02	0.41
1:B:195:ARG:HG2	1:B:225:LEU:CD1	2.50	0.41
1:A:134:VAL:HG22	1:A:142:LEU:HD11	2.03	0.41
1:B:93:TYR:CG	1:B:125:LEU:HD21	2.55	0.41
1:B:142:LEU:HD13	1:B:172:ILE:HD13	2.02	0.41
1:A:54:ARG:HH22	1:A:95:ASP:CG	2.29	0.41
1:B:110:TRP:CH2	1:B:159:VAL:HG21	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:ASN:HD21	1:B:228:ASN:HD22	1.67	0.41
1:B:211:SER:HB2	1:B:213:ILE:HD12	2.02	0.41
1:A:161:ASN:HD22	1:A:161:ASN:HA	1.50	0.40
1:A:94:GLU:O	1:A:98:LYS:HG2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	235/249 (94%)	228 (97%)	7 (3%)	0	100	100
1	B	244/249 (98%)	239 (98%)	5 (2%)	0	100	100
All	All	479/498 (96%)	467 (98%)	12 (2%)	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	201/221 (91%)	185 (92%)	16 (8%)	11	13
1	B	214/221 (97%)	196 (92%)	18 (8%)	10	11
All	All	415/442 (94%)	381 (92%)	34 (8%)	10	12

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	PRO
1	A	10	GLU
1	A	21	SER
1	A	22	LEU
1	A	45	ASN
1	A	96	LEU
1	A	97	ARG
1	A	108	LEU
1	A	110	TRP
1	A	125	LEU
1	A	136	ILE
1	A	142	LEU
1	A	160	ILE
1	A	161	ASN
1	A	214	SER
1	A	247	ILE
1	B	10	GLU
1	B	22	LEU
1	B	28	ARG
1	B	42	LYS
1	B	45	ASN
1	B	60	LEU
1	B	64	ARG
1	B	98	LYS
1	B	110	TRP
1	B	125	LEU
1	B	137	LEU
1	B	142	LEU
1	B	158	ILE
1	B	160	ILE
1	B	170	LEU
1	B	192	GLU
1	B	215	GLU
1	B	247	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	118	GLN
1	A	161	ASN

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Mol	Chain	Res	Type
1	B	41	ASN
1	B	118	GLN
1	B	124	ASN
1	B	161	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	B	301	-	4,4,4	1.17	0	6,6,6	1.37	1 (16%)
2	PO4	A	300	-	4,4,4	0.46	0	6,6,6	1.75	2 (33%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300	PO4	O3-P-O1	-3.27	99.40	110.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300	PO4	O4-P-O2	2.35	115.22	107.91
2	B	301	PO4	O4-P-O1	-2.01	103.83	110.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	PO4	1	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	241/249 (96%)	-0.07	9 (3%) 45 42	11, 21, 39, 50	0
1	B	246/249 (98%)	-0.08	10 (4%) 41 38	11, 22, 43, 56	0
All	All	487/498 (97%)	-0.07	19 (3%) 43 40	11, 22, 42, 56	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	211	SER	4.0
1	A	211	SER	3.7
1	A	24	ALA	3.2
1	B	61	ASN	3.0
1	B	23	HIS	2.9
1	B	25	SER	2.8
1	A	164	ASN	2.6
1	B	193	ASN	2.6
1	A	23	HIS	2.6
1	B	27	GLN	2.4
1	A	57	PRO	2.3
1	B	97	ARG	2.3
1	A	193	ASN	2.2
1	B	22	LEU	2.2
1	A	184	LEU	2.1
1	B	221	GLU	2.1
1	A	247	ILE	2.1
1	A	22	LEU	2.0
1	B	28	ARG	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 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($\text{\AA}^2$)	Q<0.9
2	PO4	A	300	5/5	0.94	0.08	34,35,41,41	0
2	PO4	B	301	5/5	0.94	0.08	32,37,39,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.