



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:21 AM UTC

PDB ID : 2SFP / pdb_00002sfp
Title : ALANINE RACEMASE WITH BOUND PROPIONATE INHIBITOR
Authors : Morollo, A.A.; Petsko, G.A.; Ringe, D.
Deposited on : 1999-02-16
Resolution : 1.90 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

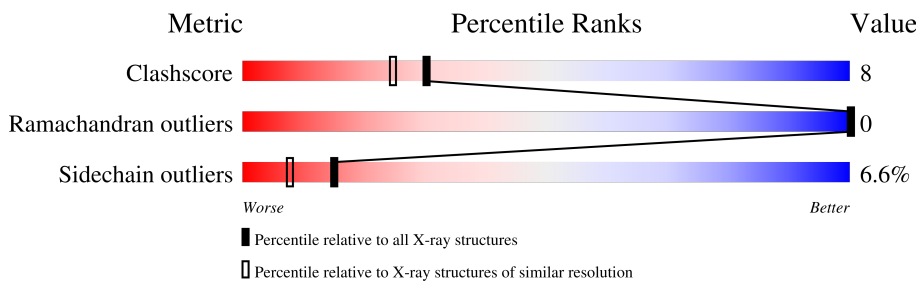
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	388	 75% 20% . .
1	B	388	 79% 16% . .

2 Entry composition [i](#)

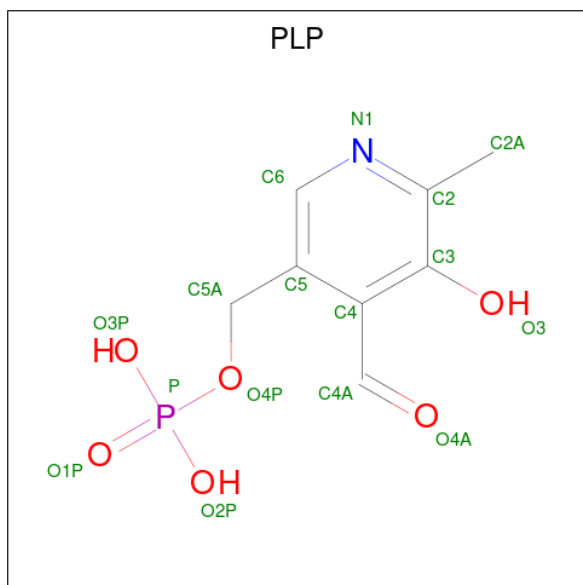
There are 4 unique types of molecules in this entry. The entry contains 6279 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (ALANINE RACEMASE).

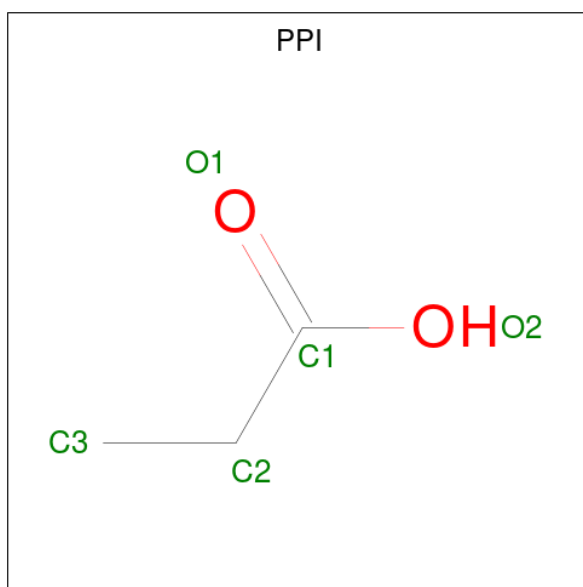
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			3016	1932	534	537	13			
1	B	379	Total	C	N	O	S	0	0	0
			3016	1932	534	537	13			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is PROPANOIC ACID (CCD ID: PPI) (formula: $C_3H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			5	3	2		
3	B	1	Total	C	O	0	0
			5	3	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	102	Total	O	0	0
			102	102		
4	B	105	Total	O	0	0
			105	105		

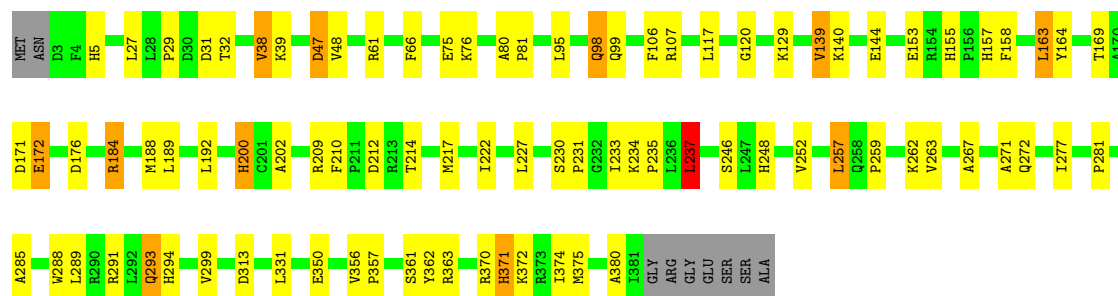
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

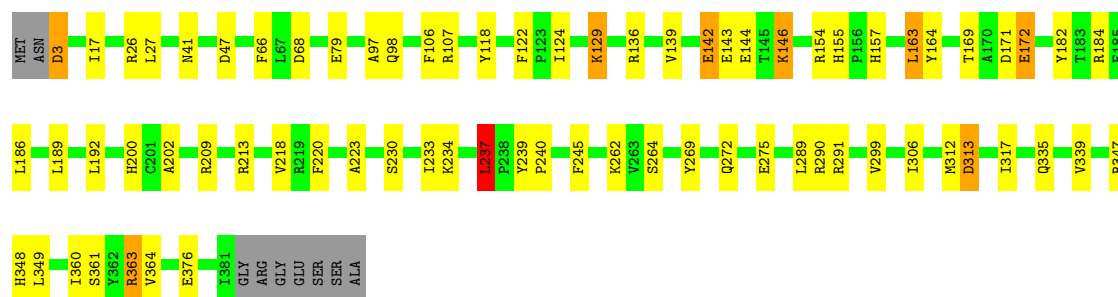
• Molecule 1: PROTEIN (ALANINE RACEMASE)

Chain A: 



• Molecule 1: PROTEIN (ALANINE RACEMASE)

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.70Å 90.00Å 85.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.90	Depositor
% Data completeness (in resolution range)	82.9 (10.00-1.90)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.205 , 0.259	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6279	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: KCX, PPI, PLP

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	0.44	0/3081	0.95	14/4188 (0.3%)
1	B	0.45	0/3081	0.94	11/4188 (0.3%)
All	All	0.44	0/6162	0.95	25/8376 (0.3%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	361	SER	N-CA-C	9.24	121.78	110.41
1	B	361	SER	N-CA-C	9.10	121.61	110.41
1	B	142	GLU	N-CA-C	7.38	118.97	111.07
1	B	47	ASP	N-CA-C	7.35	119.99	111.02
1	B	360	ILE	N-CA-C	-6.77	99.05	108.12
1	B	237	LEU	CA-C-N	6.66	126.44	119.05
1	B	237	LEU	C-N-CA	6.66	126.44	119.05
1	A	47	ASP	N-CA-C	6.55	119.02	111.02
1	B	289	LEU	N-CA-C	6.13	118.21	110.24
1	A	202	ALA	N-CA-C	6.12	119.88	110.20
1	A	237	LEU	CA-C-N	6.03	125.91	119.28
1	A	237	LEU	C-N-CA	6.03	125.91	119.28
1	A	169	THR	N-CA-C	5.83	118.34	110.88
1	B	245	PHE	N-CA-C	5.68	118.50	109.24
1	A	371	HIS	N-CA-C	-5.59	105.76	112.58
1	B	364	VAL	N-CA-C	5.51	113.32	107.76
1	B	202	ALA	N-CA-C	5.44	118.79	110.20
1	B	234	LYS	N-CA-C	5.39	120.26	113.25
1	A	263	VAL	N-CA-C	5.31	115.77	108.12
1	A	289	LEU	N-CA-C	5.28	117.10	110.24
1	A	176	ASP	N-CA-C	5.14	116.57	111.07
1	A	285	ALA	N-CA-C	-5.09	106.90	113.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	293	GLN	N-CA-C	-5.07	102.87	110.23
1	A	5	HIS	N-CA-C	5.06	119.00	111.87
1	A	99	GLN	N-CA-C	-5.05	106.71	112.92

There are no chirality outliers.

There are no planarity outliers.

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	3016	0	3003	62	0
1	B	3016	0	3004	46	0
2	A	15	0	7	0	0
2	B	15	0	7	0	0
3	A	5	0	5	0	0
3	B	5	0	5	1	0
4	A	102	0	0	0	0
4	B	105	0	0	1	0
All	All	6279	0	6031	101	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 8.

All (101) 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:231:PRO:HA	1:A:234:LYS:HE3	1.50	0.90
1:B:169:THR:HB	1:B:172:GLU:HG3	1.55	0.87
1:B:163:LEU:HD13	1:B:192:LEU:HD11	1.57	0.86
1:A:163:LEU:HD13	1:A:192:LEU:HD11	1.61	0.81
1:B:129:KCX:OQ1	1:B:136:ARG:HB3	1.83	0.78
1:A:48:VAL:HG11	1:A:76:LYS:CB	2.15	0.76
1:A:155:HIS:HD2	1:A:157:HIS:H	1.33	0.76
1:B:97:ALA:HB2	1:B:124:ILE:HG12	1.68	0.75
1:B:142:GLU:CG	1:B:146:LYS:HE3	2.19	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:VAL:O	1:A:38:VAL:HG13	1.91	0.71
1:A:48:VAL:HG11	1:A:76:LYS:HB3	1.74	0.70
1:A:61:ARG:HG2	1:A:81:PRO:HG2	1.73	0.70
1:A:230:SER:O	1:A:233:ILE:HG22	1.92	0.69
1:B:171:ASP:HB2	1:B:230:SER:HB3	1.75	0.67
1:B:155:HIS:HD2	1:B:157:HIS:H	1.42	0.67
1:B:163:LEU:HD11	1:B:189:LEU:CD2	2.26	0.66
1:A:246:SER:OG	1:A:248:HIS:HE1	1.78	0.65
1:A:293:GLN:O	1:A:294:HIS:HB2	1.96	0.65
1:A:231:PRO:HA	1:A:234:LYS:CE	2.25	0.64
1:A:38:VAL:O	1:A:38:VAL:CG1	2.46	0.62
1:A:163:LEU:HD11	1:A:189:LEU:HD23	1.81	0.62
1:A:350:GLU:HB2	1:B:291:ARG:NH2	2.16	0.60
1:B:98:GLN:HG3	1:B:122:PHE:CE2	2.38	0.59
1:B:163:LEU:HD11	1:B:189:LEU:HD23	1.86	0.57
1:A:262:LYS:HD3	1:A:267:ALA:HB1	1.86	0.57
1:A:184:ARG:O	1:A:188:MET:HG3	2.03	0.57
1:A:39:LYS:HE3	1:B:313:ASP:OD2	2.04	0.56
1:A:48:VAL:HG11	1:A:76:LYS:HB2	1.87	0.56
1:A:299:VAL:O	1:A:299:VAL:HG13	2.04	0.56
1:B:142:GLU:HG2	1:B:146:LYS:HE3	1.86	0.56
1:A:291:ARG:NH2	1:B:349:LEU:O	2.40	0.55
1:A:48:VAL:CG1	1:A:76:LYS:HD3	2.36	0.55
1:A:350:GLU:HB2	1:B:291:ARG:HH22	1.71	0.55
1:B:209:ARG:HG3	1:B:233:ILE:HD11	1.90	0.54
1:A:293:GLN:O	1:A:294:HIS:CB	2.56	0.54
1:A:155:HIS:CD2	1:A:157:HIS:H	2.21	0.54
1:A:234:LYS:HA	1:A:237:LEU:HD22	1.89	0.54
1:B:291:ARG:HB2	1:B:291:ARG:HH11	1.73	0.53
1:A:66:PHE:HZ	1:B:363:ARG:HG2	1.74	0.53
1:A:246:SER:OG	1:A:248:HIS:CE1	2.61	0.53
1:B:291:ARG:HB2	1:B:291:ARG:NH1	2.24	0.52
1:A:288:TRP:CE2	1:A:331:LEU:HD23	2.45	0.52
1:A:370:ARG:O	1:A:371:HIS:HB2	2.10	0.51
1:B:299:VAL:O	1:B:299:VAL:HG13	2.11	0.51
1:A:233:ILE:O	1:A:237:LEU:HD13	2.13	0.49
1:B:26:ARG:HH11	1:B:26:ARG:HG3	1.77	0.49
1:A:209:ARG:HG3	1:A:233:ILE:HD11	1.95	0.49
1:B:142:GLU:HG3	1:B:146:LYS:HE3	1.93	0.49
1:A:29:PRO:HG2	1:A:32:THR:OG1	2.13	0.48
1:A:80:ALA:HB1	1:A:81:PRO:HD2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:TYR:OH	1:B:122:PHE:O	2.20	0.48
1:B:312:MET:HE3	3:B:400:PPI:H32	1.95	0.48
1:A:29:PRO:HB2	1:A:31:ASP:OD2	2.14	0.47
1:A:106:PHE:CD2	1:A:107:ARG:HG3	2.50	0.47
1:A:363:ARG:HG2	1:B:66:PHE:HZ	1.80	0.47
1:A:200:HIS:HB3	1:A:217:MET:HB3	1.96	0.47
1:B:299:VAL:O	1:B:299:VAL:CG1	2.62	0.47
1:B:220:PHE:CE1	1:B:223:ALA:HB3	2.50	0.47
1:B:239:TYR:HB2	1:B:240:PRO:HD2	1.97	0.47
1:B:142:GLU:HG2	1:B:146:LYS:CE	2.43	0.47
1:A:98:GLN:HE21	1:A:98:GLN:HB3	1.49	0.46
1:A:231:PRO:CA	1:A:234:LYS:HE3	2.34	0.46
1:A:164:TYR:HB2	1:A:200:HIS:O	2.15	0.46
1:B:142:GLU:HG2	1:B:146:LYS:NZ	2.30	0.46
1:A:163:LEU:HD11	1:A:189:LEU:CD2	2.44	0.45
1:A:231:PRO:HB3	1:A:234:LYS:NZ	2.31	0.45
1:B:230:SER:O	1:B:233:ILE:HG22	2.16	0.45
1:A:48:VAL:HG13	1:A:76:LYS:HD3	1.98	0.45
1:B:139:VAL:HG23	1:B:144:GLU:HG2	1.98	0.45
1:A:331:LEU:HD12	1:A:331:LEU:N	2.31	0.45
1:A:120:GLY:HA3	1:A:157:HIS:CE1	2.52	0.44
1:A:362:TYR:OH	1:B:41:ASN:HB2	2.17	0.44
1:A:222:ILE:HG13	1:A:227:LEU:HB2	2.00	0.44
1:B:306:ILE:HA	1:B:317:ILE:HG22	1.98	0.44
1:A:234:LYS:N	1:A:235:PRO:HD2	2.32	0.44
1:A:252:VAL:HG11	1:A:281:PRO:HG3	2.00	0.43
1:B:182:TYR:CE2	1:B:186:LEU:HD11	2.53	0.43
1:A:288:TRP:CZ2	1:A:331:LEU:HD23	2.54	0.43
1:A:200:HIS:CD2	1:A:200:HIS:N	2.87	0.43
1:A:362:TYR:O	1:A:380:ALA:HB3	2.18	0.43
1:B:106:PHE:CD2	1:B:107:ARG:HG3	2.54	0.42
1:B:262:LYS:HA	1:B:269:TYR:O	2.19	0.42
1:B:3:ASP:OD1	1:B:3:ASP:N	2.51	0.42
1:A:139:VAL:CG2	1:A:144:GLU:HB3	2.50	0.42
1:A:155:HIS:HB3	1:A:158:PHE:HB2	2.00	0.42
1:B:237:LEU:HD12	1:B:237:LEU:HA	1.93	0.42
1:B:347:ARG:HH21	1:B:348:HIS:CE1	2.37	0.42
1:A:257:LEU:HG	1:A:277:ILE:HG13	2.01	0.42
1:B:347:ARG:HH21	1:B:348:HIS:HE1	1.68	0.42
1:B:66:PHE:CE2	1:B:68:ASP:HB2	2.55	0.41
1:A:259:PRO:HA	1:A:271:ALA:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:SER:HB3	4:B:569:HOH:O	2.20	0.41
1:A:48:VAL:CG1	1:A:76:LYS:HB2	2.51	0.41
1:A:356:VAL:HB	1:A:357:PRO:HD3	2.01	0.41
1:B:79:GLU:HA	1:B:79:GLU:OE1	2.19	0.41
1:A:171:ASP:HA	1:A:233:ILE:HD12	2.03	0.41
1:A:172:GLU:O	1:A:209:ARG:NH2	2.50	0.41
1:B:164:TYR:HB2	1:B:200:HIS:O	2.21	0.41
1:B:142:GLU:CG	1:B:146:LYS:CE	2.96	0.40
1:A:47:ASP:OD1	1:A:48:VAL:N	2.54	0.40
1:A:210:PHE:O	1:A:214:THR:HG23	2.21	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	336/388 (87%)	326 (97%)	10 (3%)	0	100	100
1	B	376/388 (97%)	366 (97%)	10 (3%)	0	100	100
All	All	712/776 (92%)	692 (97%)	20 (3%)	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	309/322 (96%)	288 (93%)	21 (7%)	14	7
1	B	309/322 (96%)	289 (94%)	20 (6%)	15	8
All	All	618/644 (96%)	577 (93%)	41 (7%)	15	8

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	38	VAL
1	A	75	GLU
1	A	95	LEU
1	A	98	GLN
1	A	117	LEU
1	A	139	VAL
1	A	140	LYS
1	A	153	GLU
1	A	163	LEU
1	A	172	GLU
1	A	184	ARG
1	A	200	HIS
1	A	212	ASP
1	A	237	LEU
1	A	257	LEU
1	A	272	GLN
1	A	313	ASP
1	A	372	LYS
1	A	374	ILE
1	A	375	MET
1	B	3	ASP
1	B	17	ILE
1	B	27	LEU
1	B	143	GLU
1	B	146	LYS
1	B	154	ARG
1	B	163	LEU
1	B	172	GLU
1	B	184	ARG
1	B	213	ARG
1	B	218	VAL
1	B	237	LEU
1	B	272	GLN
1	B	275	GLU

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Mol	Chain	Res	Type
1	B	290	ARG
1	B	313	ASP
1	B	335	GLN
1	B	339	VAL
1	B	363	ARG
1	B	376	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	HIS
1	A	33	HIS
1	A	98	GLN
1	A	155	HIS
1	A	248	HIS
1	A	348	HIS
1	B	5	HIS
1	B	155	HIS
1	B	248	HIS
1	B	258	GLN
1	B	314	GLN
1	B	348	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 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
1	KCX	A	129	1	10,11,12	0.89	0	6,12,14	2.14	1 (16%)
1	KCX	B	129	1	10,11,12	1.31	1 (10%)	6,12,14	1.91	1 (16%)

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
1	KCX	A	129	1	-	1/9/10/12	-
1	KCX	B	129	1	-	2/9/10/12	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	129	KCX	CX-NZ	-2.71	1.30	1.35

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	KCX	OQ1-CX-NZ	-4.83	117.58	124.92
1	B	129	KCX	OQ1-CX-NZ	-4.08	118.72	124.92

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	129	KCX	CE-CD-CG-CB
1	B	129	KCX	CG-CD-CE-NZ
1	B	129	KCX	CE-CD-CG-CB

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	129	KCX	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	PLP	A	500	1	15,15,16	1.40	1 (6%)	21,22,23	1.97	6 (28%)
2	PLP	B	500	1	15,15,16	1.93	6 (40%)	21,22,23	1.84	4 (19%)
3	PPI	B	400	-	4,4,4	0.98	0	4,4,4	0.75	0
3	PPI	A	400	-	4,4,4	1.01	0	4,4,4	0.75	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	PLP	A	500	1	-	2/6/6/8	0/1/1/1
2	PLP	B	500	1	-	2/6/6/8	0/1/1/1
3	PPI	B	400	-	-	2/2/2/2	-
3	PPI	A	400	-	-	1/2/2/2	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	PLP	C2A-C2	3.96	1.56	1.50
2	B	500	PLP	C5A-C5	-2.97	1.43	1.50
2	A	500	PLP	C5-C4	-2.88	1.37	1.40
2	B	500	PLP	C6-N1	2.40	1.39	1.34
2	B	500	PLP	O3-C3	2.35	1.42	1.36
2	B	500	PLP	P-O2P	-2.31	1.46	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	PLP	P-O3P	-2.04	1.47	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	PLP	C4A-C4-C5	5.73	126.84	120.94
2	A	500	PLP	C6-C5-C4	5.18	122.35	118.10
2	A	500	PLP	C5A-C5-C6	-4.24	112.46	119.36
2	B	500	PLP	C4A-C4-C3	-3.58	114.56	120.52
2	A	500	PLP	C4A-C4-C5	2.76	123.78	120.94
2	B	500	PLP	O3P-P-O2P	2.48	117.11	107.80
2	B	500	PLP	O2P-P-O4P	-2.38	100.47	106.67
2	A	500	PLP	O2P-P-O4P	-2.33	100.60	106.67
2	A	500	PLP	C3-C4-C5	-2.14	116.02	118.59
2	A	500	PLP	O3-C3-C4	2.02	123.38	118.10

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	500	PLP	C6-C5-C5A-O4P
2	A	500	PLP	C4-C5-C5A-O4P
2	B	500	PLP	C4-C5-C5A-O4P
3	A	400	PPI	O2-C1-C2-C3
3	B	400	PPI	O2-C1-C2-C3
2	A	500	PLP	C6-C5-C5A-O4P
3	B	400	PPI	O1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	400	PPI	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.