



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:27 AM UTC

PDB ID : 7DFB / pdb_00007dfb
Title : Crystal of Arrestin2-V2Rpp-6-7-Fab30 complex
Authors : Sun, J.P.; Yu, X.; Xiao, P.; He, Q.T.; Lin, J.Y.; Zhu, Z.L.
Deposited on : 2020-11-06
Resolution : 3.28 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)
Xtrriage (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

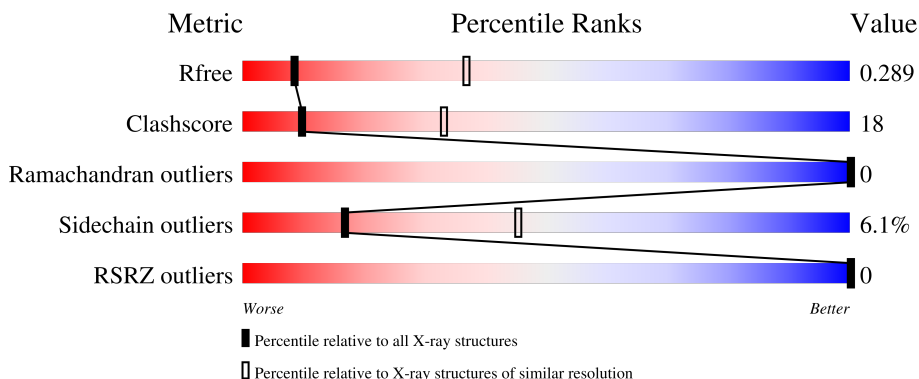
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.28 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	1303 (3.30-3.26)
Clashscore	190562	1354 (3.30-3.26)
Ramachandran outliers	187476	1334 (3.30-3.26)
Sidechain outliers	187428	1333 (3.30-3.26)
RSRZ outliers	180081	1303 (3.30-3.26)

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	426	 51% 31% 16%
2	V	23	 43% 30% 13% 13%
3	L	227	 56% 33% 7%
4	H	249	 57% 26% 14%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5913 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 Beta-arrestin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	356	2702	1728	462	502	10	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	419	LEU	-	expression tag	UNP P17870
A	420	GLU	-	expression tag	UNP P17870
A	421	HIS	-	expression tag	UNP P17870
A	422	HIS	-	expression tag	UNP P17870
A	423	HIS	-	expression tag	UNP P17870
A	424	HIS	-	expression tag	UNP P17870
A	425	HIS	-	expression tag	UNP P17870
A	426	HIS	-	expression tag	UNP P17870

- Molecule 2 is a protein called V2Rpp-6-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	P				S
2	V	20	151	75	21	48	6	1	0	0	0

- Molecule 3 is a protein called FAB30 LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	L	212	1550	970	254	321	5	0	0	0

- Molecule 4 is a protein called FAB30 HEAVY CHAIN.

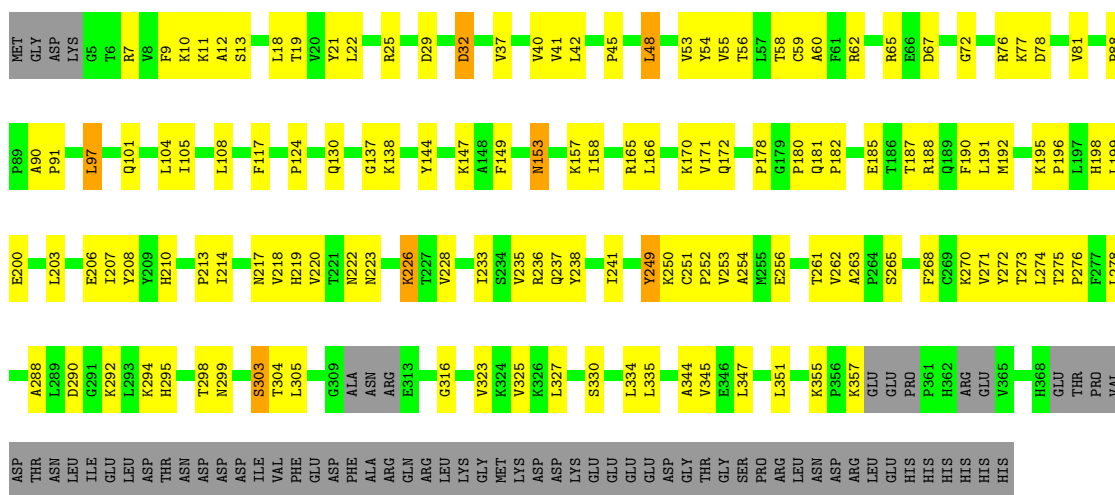
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	H	213	1510	957	253	295	5	0	0	0

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.

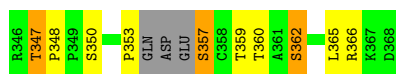
- Molecule 1: Beta-arrestin-1

Chain A: 



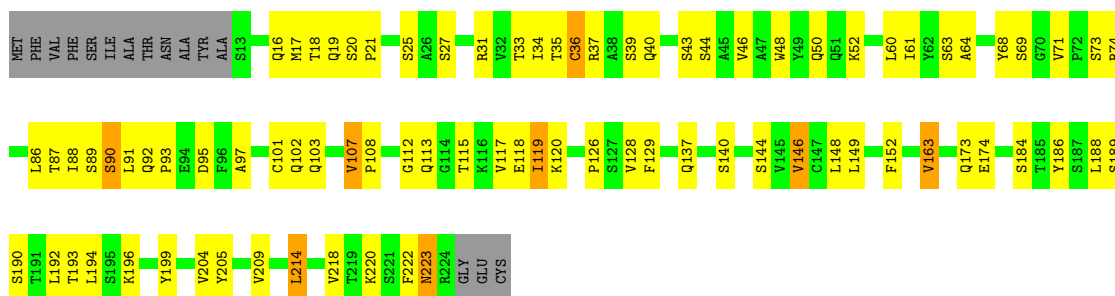
- Molecule 2: V2Rpp-6-7

Chain V: 



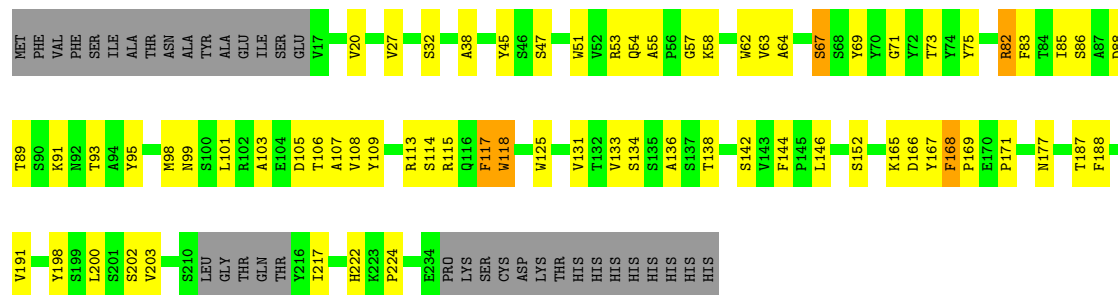
- Molecule 3: FAB30 LIGHT CHAIN

Chain L: 



- Molecule 4: FAB30 HEAVY CHAIN

Chain H:  57% 26% 14%



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	115.48Å 116.88Å 143.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.36 – 3.28 45.36 – 3.28	Depositor EDS
% Data completeness (in resolution range)	98.5 (45.36-3.28) 99.1 (45.36-3.28)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 3.25Å)	Xtrriage
Refinement program	PHENIX 1.14 3260	Depositor
R, R_{free}	0.231 , 0.290 0.234 , 0.289	Depositor DCC
R_{free} test set	750 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	110.1	Xtrriage
Anisotropy	0.505	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 83.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.008 for -k,-h,-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5913	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.48	0/2757	0.72	0/3751
2	V	0.46	0/85	0.73	0/107
3	L	0.44	0/1581	0.68	0/2158
4	H	0.45	0/1550	0.77	4/2124 (0.2%)
All	All	0.46	0/5973	0.72	4/8140 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	118	TRP	CA-C-N	6.19	133.36	121.54
4	H	118	TRP	C-N-CA	6.19	133.36	121.54
4	H	117	PHE	CB-CA-C	-5.98	100.92	110.14
4	H	168	PHE	N-CA-C	5.07	121.02	109.81

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	2702	0	2697	95	0
2	V	151	0	104	6	0
3	L	1550	0	1463	59	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	1510	0	1353	55	0
All	All	5913	0	5617	203	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 18.

All (203) 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)
4:H:82:ARG:HG2	4:H:82:ARG:HH11	1.40	0.85
4:H:134:SER:HB3	4:H:168:PHE:HZ	1.42	0.84
1:A:62:ARG:HG2	1:A:76:ARG:HG2	1.69	0.74
1:A:40:VAL:HG22	1:A:104:LEU:HD13	1.70	0.73
3:L:18:THR:HG22	3:L:113:GLN:HE22	1.54	0.73
1:A:42:LEU:HD22	1:A:108:LEU:HD13	1.71	0.72
3:L:90:SER:O	3:L:92:GLN:NE2	2.22	0.72
1:A:292:LYS:HG2	1:A:298:THR:OG1	1.88	0.72
3:L:146:VAL:HG11	4:H:146:LEU:HD13	1.73	0.71
4:H:134:SER:HB3	4:H:168:PHE:CZ	2.26	0.69
4:H:53:ARG:HB2	4:H:109:TYR:CE1	2.29	0.67
1:A:233:ILE:HD13	1:A:325:VAL:HG22	1.76	0.66
3:L:68:TYR:HB3	3:L:71:VAL:HG21	1.78	0.66
3:L:214:LEU:HD22	3:L:218:VAL:HG21	1.78	0.66
3:L:48:TRP:HB2	3:L:61:ILE:HG22	1.77	0.66
3:L:128:VAL:HG21	3:L:209:VAL:HG21	1.76	0.65
4:H:82:ARG:HH11	4:H:82:ARG:CG	2.10	0.65
1:A:213:PRO:HA	1:A:275:THR:HG22	1.78	0.65
1:A:12:ALA:HA	1:A:19:THR:HA	1.77	0.65
3:L:34:ILE:HD12	3:L:86:LEU:HD23	1.78	0.64
1:A:171:VAL:HG22	1:A:172:GLN:H	1.62	0.63
1:A:29:ASP:HB2	1:A:170:LYS:HE2	1.80	0.63
1:A:355:LYS:HD2	1:A:357:LYS:HE3	1.80	0.63
3:L:204:VAL:HG13	3:L:204:VAL:O	1.98	0.62
1:A:203:LEU:HD21	1:A:323:VAL:HG11	1.81	0.62
4:H:55:ALA:HB3	4:H:58:LYS:HB2	1.80	0.62
1:A:222:ASN:O	1:A:265:SER:HA	2.00	0.61
4:H:98:MET:HE1	4:H:131:VAL:HG11	1.83	0.60
3:L:31:ARG:HG2	3:L:89:SER:HA	1.82	0.60
4:H:103:ALA:O	4:H:106:THR:HG22	2.02	0.59
1:A:182:PRO:HG2	1:A:203:LEU:HB2	1.84	0.59
1:A:217:ASN:HA	1:A:271:VAL:HG12	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:165:LYS:NZ	4:H:166:ASP:OD2	2.26	0.58
1:A:185:GLU:HB2	1:A:198:HIS:HE1	1.68	0.58
3:L:33:THR:HG22	3:L:87:THR:HA	1.86	0.58
3:L:107:VAL:N	3:L:108:PRO:HD2	2.19	0.58
1:A:235:VAL:HG11	1:A:276:PRO:HG3	1.85	0.57
4:H:142:SER:HB3	4:H:144:PHE:CZ	2.39	0.57
1:A:88:PRO:HG2	1:A:91:PRO:HD3	1.86	0.57
3:L:199:TYR:HH	3:L:222:PHE:HE2	1.51	0.57
3:L:92:GLN:HB3	3:L:93:PRO:HD2	1.85	0.57
1:A:199:LEU:HD13	1:A:327:LEU:HD11	1.87	0.56
4:H:75:TYR:OH	4:H:85:ILE:N	2.33	0.56
3:L:18:THR:O	3:L:36:CYS:HA	2.06	0.55
4:H:53:ARG:NH1	4:H:105:ASP:HA	2.22	0.55
3:L:222:PHE:CD2	3:L:222:PHE:C	2.85	0.55
3:L:204:VAL:HB	3:L:223:ASN:HD21	1.71	0.55
4:H:118:TRP:CD1	4:H:118:TRP:N	2.71	0.55
1:A:13:SER:OG	1:A:18:LEU:HB2	2.07	0.55
4:H:67:SER:O	4:H:71:GLY:N	2.39	0.55
1:A:54:TYR:HB2	1:A:149:PHE:CZ	2.43	0.55
3:L:27:SER:HB3	3:L:120:LYS:HE2	1.89	0.54
3:L:16:GLN:H	3:L:39:SER:HB2	1.73	0.54
1:A:299:ASN:HB3	4:H:117:PHE:HE1	1.72	0.54
1:A:200:GLU:HB2	1:A:219:HIS:HB3	1.89	0.54
3:L:118:GLU:HG3	3:L:186:TYR:OH	2.08	0.54
3:L:46:VAL:HG13	3:L:102:GLN:O	2.08	0.54
3:L:50:GLN:HB3	3:L:60:LEU:HD11	1.89	0.54
1:A:77:LYS:NZ	2:V:347:TPO:HG23	2.23	0.53
1:A:210:HIS:HB3	4:H:117:PHE:CZ	2.44	0.53
1:A:18:LEU:HD22	1:A:41:VAL:CG2	2.38	0.53
3:L:19:GLN:HA	3:L:35:THR:O	2.09	0.53
1:A:124:PRO:HG3	1:A:316:GLY:HA3	1.89	0.53
4:H:188:PHE:O	4:H:200:LEU:HD11	2.09	0.53
1:A:78:ASP:OD1	1:A:78:ASP:N	2.42	0.52
3:L:119:ILE:HD12	3:L:184:SER:OG	2.10	0.52
1:A:237:GLN:N	1:A:251:CYS:O	2.42	0.52
3:L:126:PRO:HB3	3:L:152:PHE:CD2	2.43	0.52
1:A:222:ASN:ND2	1:A:263:ALA:O	2.43	0.52
3:L:20:SER:HB2	3:L:35:THR:HG23	1.91	0.52
4:H:82:ARG:HG2	4:H:82:ARG:NH1	2.19	0.52
3:L:149:LEU:HB2	3:L:188:LEU:HB3	1.92	0.52
4:H:138:THR:HG23	4:H:169:PRO:HG3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:LYS:O	1:A:158:ILE:HD13	2.11	0.51
4:H:88:ASP:HB2	4:H:95:TYR:HE2	1.76	0.51
1:A:153:ASN:N	1:A:153:ASN:OD1	2.43	0.51
1:A:295:HIS:NE2	2:V:353:PRO:HG3	2.25	0.51
1:A:149:PHE:CD2	1:A:158:ILE:HD12	2.46	0.50
1:A:45:PRO:HA	1:A:48:LEU:HB3	1.94	0.50
1:A:237:GLN:OE1	1:A:288:ALA:HB3	2.12	0.50
3:L:204:VAL:CB	3:L:223:ASN:HD21	2.25	0.50
4:H:115:ARG:C	4:H:117:PHE:H	2.19	0.50
1:A:187:THR:HG23	1:A:198:HIS:HA	1.93	0.50
1:A:238:TYR:CE1	1:A:250:LYS:HB3	2.47	0.50
1:A:188:ARG:HD2	1:A:190:PHE:CE1	2.46	0.50
1:A:196:PRO:O	1:A:223:ASN:HB2	2.11	0.50
4:H:83:PHE:CE1	4:H:98:MET:HB3	2.47	0.50
1:A:72:GLY:HA2	1:A:137:GLY:HA3	1.94	0.49
1:A:181:GLN:HG3	1:A:203:LEU:O	2.11	0.49
1:A:206:GLU:O	1:A:206:GLU:HG2	2.12	0.49
3:L:148:LEU:HD11	4:H:203:VAL:HG21	1.93	0.49
1:A:236:ARG:HA	1:A:252:PRO:HA	1.94	0.49
4:H:136:ALA:HB3	4:H:168:PHE:CE2	2.47	0.49
4:H:187:THR:HG23	4:H:202:SER:HB2	1.95	0.48
1:A:60:ALA:HA	1:A:78:ASP:HA	1.95	0.48
4:H:62:TRP:HE1	4:H:64:ALA:C	2.22	0.47
1:A:65:ARG:HG2	1:A:67:ASP:OD1	2.13	0.47
1:A:236:ARG:NH1	1:A:250:LYS:HE3	2.28	0.47
3:L:174:GLU:HA	3:L:190:SER:HA	1.96	0.47
1:A:97:LEU:HB2	1:A:101:GLN:HB2	1.95	0.47
2:V:362:SEP:O1P	3:L:43:SER:OG	2.30	0.47
3:L:17:MET:HA	3:L:37:ARG:O	2.14	0.47
1:A:191:LEU:HD23	1:A:334:LEU:CD2	2.44	0.47
1:A:305:LEU:HD23	1:A:305:LEU:HA	1.74	0.47
4:H:177:ASN:N	4:H:217:ILE:O	2.40	0.47
1:A:37:VAL:HG13	1:A:117:PHE:HB2	1.97	0.46
3:L:74:ARG:HD2	3:L:88:ILE:HD11	1.96	0.46
4:H:222:HIS:CD2	4:H:224:PRO:HD2	2.50	0.46
1:A:233:ILE:CD1	1:A:325:VAL:HG22	2.42	0.46
3:L:48:TRP:CZ3	3:L:101:CYS:HB3	2.51	0.46
3:L:118:GLU:HG2	3:L:119:ILE:H	1.80	0.46
1:A:10:LYS:HB3	1:A:21:TYR:CD2	2.51	0.46
3:L:188:LEU:HD23	3:L:189:SER:N	2.30	0.46
4:H:167:TYR:O	4:H:198:TYR:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:20:VAL:HG22	4:H:38:ALA:HB3	1.97	0.46
1:A:55:VAL:HG12	1:A:117:PHE:HE1	1.81	0.46
1:A:11:LYS:HD2	1:A:166:LEU:HD13	1.97	0.46
1:A:325:VAL:O	1:A:344:ALA:HA	2.15	0.46
1:A:10:LYS:HB3	1:A:21:TYR:CE2	2.50	0.46
3:L:220:LYS:HE3	3:L:220:LYS:HB3	1.78	0.46
1:A:7:ARG:HA	1:A:7:ARG:HD2	1.79	0.46
1:A:59:CYS:HB2	1:A:144:TYR:CE1	2.51	0.45
1:A:323:VAL:HG12	1:A:347:LEU:HB2	1.98	0.45
3:L:17:MET:O	3:L:112:GLY:HA2	2.16	0.45
1:A:254:ALA:HB2	1:A:276:PRO:HA	1.98	0.45
4:H:113:ARG:HG2	4:H:114:SER:N	2.32	0.45
3:L:52:LYS:HG2	3:L:97:ALA:HB2	1.97	0.45
1:A:208:TYR:O	1:A:351:LEU:HD12	2.17	0.44
3:L:126:PRO:O	3:L:128:VAL:HG23	2.16	0.44
4:H:47:SER:HB3	4:H:113:ARG:HG3	1.99	0.44
3:L:163:VAL:HA	3:L:205:TYR:HA	1.98	0.44
1:A:290:ASP:C	1:A:298:THR:HG21	2.43	0.44
4:H:57:GLY:C	4:H:58:LYS:HD3	2.42	0.44
4:H:109:TYR:CD2	4:H:131:VAL:HB	2.52	0.44
4:H:82:ARG:CG	4:H:82:ARG:NH1	2.72	0.44
4:H:125:TRP:CD1	4:H:125:TRP:N	2.85	0.44
1:A:60:ALA:CB	1:A:78:ASP:HA	2.48	0.44
1:A:88:PRO:HB2	1:A:90:ALA:H	1.81	0.44
3:L:91:LEU:HD23	3:L:91:LEU:H	1.82	0.44
3:L:199:TYR:OH	3:L:222:PHE:HE2	2.00	0.44
4:H:86:SER:OG	4:H:95:TYR:HB2	2.18	0.44
1:A:62:ARG:NH1	1:A:138:LYS:HG2	2.32	0.44
1:A:210:HIS:HB2	4:H:117:PHE:CE2	2.53	0.44
1:A:32:ASP:OD1	1:A:32:ASP:N	2.35	0.43
1:A:199:LEU:HD12	1:A:199:LEU:HA	1.81	0.43
1:A:58:THR:OG1	1:A:81:VAL:HG12	2.18	0.43
1:A:178:PRO:HA	1:A:207:ILE:HD12	2.00	0.43
1:A:195:LYS:HB3	1:A:223:ASN:HB3	1.99	0.43
1:A:228:VAL:HB	1:A:262:VAL:HB	2.01	0.43
3:L:63:SER:O	3:L:64:ALA:HB3	2.18	0.43
1:A:214:ILE:O	1:A:273:THR:HA	2.18	0.43
1:A:180:PRO:O	1:A:206:GLU:HB2	2.18	0.43
1:A:147:LYS:HA	1:A:165:ARG:HA	2.01	0.43
1:A:303:SER:OG	1:A:351:LEU:N	2.47	0.43
4:H:167:TYR:HE1	4:H:171:PRO:HA	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:VAL:HB	1:A:268:PHE:HB3	2.00	0.42
4:H:169:PRO:HG2	4:H:222:HIS:NE2	2.33	0.42
1:A:256:GLU:HG2	1:A:274:LEU:HD11	2.00	0.42
4:H:51:TRP:O	4:H:63:VAL:HG22	2.18	0.42
4:H:91:LYS:HD2	4:H:95:TYR:OH	2.18	0.42
3:L:102:GLN:HG2	3:L:103:GLN:N	2.35	0.42
1:A:210:HIS:CB	4:H:117:PHE:CE2	3.02	0.42
3:L:68:TYR:HB3	3:L:71:VAL:CG2	2.49	0.42
1:A:9:PHE:CD1	1:A:25:ARG:HG2	2.55	0.42
1:A:56:THR:HG22	1:A:147:LYS:HG2	2.00	0.42
1:A:294:LYS:H	1:A:294:LYS:HG2	1.74	0.42
3:L:21:PRO:HG2	3:L:34:ILE:HG23	2.01	0.42
3:L:144:SER:OG	3:L:193:THR:HG23	2.19	0.42
4:H:45:TYR:O	4:H:69:TYR:HB2	2.20	0.42
4:H:165:LYS:HG2	4:H:166:ASP:CG	2.44	0.42
3:L:20:SER:HB2	3:L:35:THR:CG2	2.48	0.42
1:A:210:HIS:CG	1:A:299:ASN:HD22	2.38	0.42
1:A:249:TYR:C	1:A:249:TYR:CD2	2.98	0.42
3:L:129:PHE:CD2	4:H:152:SER:HA	2.55	0.42
1:A:278:LEU:HD12	1:A:278:LEU:O	2.19	0.42
3:L:17:MET:HE2	3:L:103:GLN:HB3	2.01	0.42
1:A:270:LYS:HD3	1:A:272:TYR:OH	2.20	0.41
3:L:222:PHE:O	3:L:222:PHE:CG	2.70	0.41
4:H:32:SER:OG	4:H:99:ASN:HA	2.20	0.41
1:A:226:LYS:HG2	1:A:330:SER:O	2.20	0.41
1:A:199:LEU:HD23	1:A:345:VAL:HG23	2.03	0.41
1:A:241:ILE:HD12	1:A:241:ILE:N	2.35	0.41
2:V:347:TPO:HA	2:V:348:PRO:HD3	1.95	0.41
3:L:33:THR:CG2	3:L:87:THR:HG22	2.51	0.41
3:L:173:GLN:HB3	4:H:191:VAL:HG11	2.02	0.41
3:L:196:LYS:HB3	3:L:196:LYS:HE2	1.74	0.41
1:A:130:GLN:OE1	1:A:278:LEU:HD11	2.20	0.41
4:H:54:GLN:C	4:H:107:ALA:HB1	2.46	0.41
1:A:165:ARG:NH2	2:V:357:SEP:O3P	2.54	0.41
2:V:365:LEU:HG	2:V:366:ARG:H	1.86	0.41
3:L:128:VAL:CG2	3:L:209:VAL:HG21	2.47	0.41
4:H:109:TYR:CE2	4:H:131:VAL:HB	2.55	0.41
3:L:107:VAL:N	3:L:108:PRO:CD	2.84	0.41
4:H:27:VAL:HG11	4:H:101:LEU:HD12	2.03	0.41
1:A:105:ILE:H	1:A:105:ILE:HG13	1.76	0.41
1:A:323:VAL:CG1	1:A:347:LEU:HB2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:223:ASN:OD1	3:L:223:ASN:N	2.54	0.41
4:H:54:GLN:HE21	4:H:108:VAL:CG1	2.34	0.41
1:A:22:LEU:HD23	1:A:22:LEU:HA	1.97	0.40
3:L:137:GLN:O	3:L:140:SER:OG	2.34	0.40
4:H:58:LYS:HB2	4:H:58:LYS:HE2	1.91	0.40
1:A:138:LYS:HD2	1:A:138:LYS:HA	1.87	0.40
4:H:113:ARG:HG2	4:H:114:SER:O	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	348/426 (82%)	320 (92%)	28 (8%)	0	100	100
2	V	11/23 (48%)	7 (64%)	4 (36%)	0	100	100
3	L	210/227 (92%)	181 (86%)	29 (14%)	0	100	100
4	H	209/249 (84%)	185 (88%)	24 (12%)	0	100	100
All	All	778/925 (84%)	693 (89%)	85 (11%)	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	292/380 (77%)	278 (95%)	14 (5%)	23	52
2	V	9/15 (60%)	9 (100%)	0	100	100
3	L	172/199 (86%)	154 (90%)	18 (10%)	6	25
4	H	148/209 (71%)	142 (96%)	6 (4%)	27	55
All	All	621/803 (77%)	583 (94%)	38 (6%)	17	45

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ASP
1	A	48	LEU
1	A	53	VAL
1	A	97	LEU
1	A	153	ASN
1	A	192	MET
1	A	218	VAL
1	A	226	LYS
1	A	249	TYR
1	A	253	VAL
1	A	261	THR
1	A	303	SER
1	A	304	THR
1	A	335	LEU
3	L	25	SER
3	L	36	CYS
3	L	40	GLN
3	L	44	SER
3	L	69	SER
3	L	73	SER
3	L	90	SER
3	L	95	ASP
3	L	107	VAL
3	L	115	THR
3	L	117	VAL
3	L	119	ILE
3	L	146	VAL
3	L	163	VAL
3	L	192	LEU
3	L	194	LEU
3	L	214	LEU
3	L	223	ASN
4	H	67	SER

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Mol	Chain	Res	Type
4	H	73	THR
4	H	82	ARG
4	H	89	THR
4	H	93	THR
4	H	133	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	225	ASN
1	A	299	ASN
3	L	173	GLN
3	L	223	ASN
4	H	54	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

6 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	TPO	V	347	2	8,10,11	1.80	2 (25%)	10,14,16	1.92	3 (30%)
2	SEP	V	350	2	8,9,10	1.71	2 (25%)	7,12,14	1.94	1 (14%)
2	TPO	V	359	2	8,10,11	1.83	2 (25%)	10,14,16	1.71	1 (10%)
2	SEP	V	362	2	8,9,10	1.70	1 (12%)	7,12,14	1.80	2 (28%)
2	SEP	V	357	2	8,9,10	1.80	2 (25%)	7,12,14	1.15	1 (14%)
2	TPO	V	360	2	8,10,11	1.25	1 (12%)	10,14,16	2.07	2 (20%)

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	TPO	V	347	2	-	0/9/11/13	-
2	SEP	V	350	2	-	3/6/8/10	-
2	TPO	V	359	2	-	4/9/11/13	-
2	SEP	V	362	2	-	4/6/8/10	-
2	SEP	V	357	2	-	4/6/8/10	-
2	TPO	V	360	2	-	0/9/11/13	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	V	357	SEP	P-O1P	4.00	1.62	1.50
2	V	359	TPO	P-O1P	3.67	1.61	1.50
2	V	347	TPO	P-O1P	3.63	1.61	1.50
2	V	350	SEP	P-O1P	3.55	1.61	1.50
2	V	362	SEP	P-O1P	3.44	1.61	1.50
2	V	347	TPO	P-OG1	2.31	1.63	1.59
2	V	350	SEP	P-O3P	2.19	1.62	1.54
2	V	360	TPO	P-O2P	2.18	1.62	1.54
2	V	359	TPO	P-O3P	2.17	1.62	1.54
2	V	357	SEP	P-O3P	2.00	1.62	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	V	360	TPO	P-OG1-CB	-5.41	108.61	123.33
2	V	359	TPO	P-OG1-CB	-4.73	110.46	123.33
2	V	350	SEP	OG-CB-CA	4.22	112.25	108.14
2	V	347	TPO	P-OG1-CB	-4.10	112.19	123.33
2	V	362	SEP	OG-CB-CA	3.87	111.91	108.14
2	V	347	TPO	CG2-CB-CA	-3.47	106.51	113.26
2	V	362	SEP	O3P-P-OG	2.65	113.59	106.67
2	V	357	SEP	OG-CB-CA	2.56	110.64	108.14
2	V	360	TPO	CG2-CB-CA	-2.50	108.40	113.26
2	V	347	TPO	O-C-CA	-2.18	119.16	124.77

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	V	350	SEP	N-CA-CB-OG
2	V	350	SEP	CA-CB-OG-P
2	V	357	SEP	CB-OG-P-O1P
2	V	357	SEP	CB-OG-P-O2P
2	V	357	SEP	CB-OG-P-O3P
2	V	359	TPO	N-CA-CB-OG1
2	V	359	TPO	C-CA-CB-CG2
2	V	359	TPO	O-C-CA-CB
2	V	362	SEP	C-CA-CB-OG
2	V	362	SEP	CB-OG-P-O2P
2	V	362	SEP	CB-OG-P-O3P
2	V	362	SEP	CB-OG-P-O1P
2	V	359	TPO	N-CA-CB-CG2
2	V	357	SEP	CA-CB-OG-P
2	V	350	SEP	C-CA-CB-OG

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	V	347	TPO	2	0
2	V	362	SEP	1	0
2	V	357	SEP	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	356/426 (83%)	-0.28	0 100 100	69, 107, 140, 169	0
2	V	14/23 (60%)	-0.38	0 100 100	108, 133, 150, 157	0
3	L	212/227 (93%)	-0.30	0 100 100	91, 126, 149, 164	0
4	H	213/249 (85%)	-0.41	0 100 100	75, 107, 148, 166	0
All	All	795/925 (85%)	-0.32	0 100 100	69, 113, 147, 169	0

There are no RSRZ outliers to report.

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	TPO	V	347	11/12	0.51	0.11	136,142,154,157	0
2	TPO	V	359	11/12	0.79	0.07	128,144,165,172	0
2	SEP	V	350	10/11	0.83	0.07	117,125,131,132	0
2	SEP	V	357	10/11	0.85	0.06	134,138,151,152	0
2	SEP	V	362	10/11	0.91	0.06	95,102,119,120	0
2	TPO	V	360	11/12	0.95	0.05	107,115,122,123	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.