



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:00 PM GMT

PDB ID : 3VS4
Title : Crystal structure of HCK complexed with a pyrrolo-pyrimidine inhibitor 5-(4-phenoxyphenyl)-7-(tetrahydro-2H-pyran-4-yl)-7H-pyrrolo[2,3-d]pyrimidin-4-amine
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Deposited on : 2012-04-21
Resolution : 2.75 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

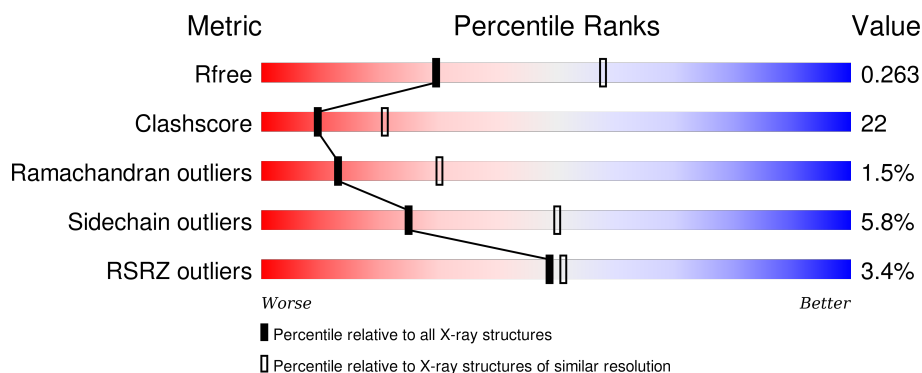
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


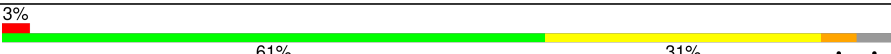
The reported resolution of this entry is 2.75 Å.

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}	91344	3050 (2.78-2.70)
Clashscore	102246	3424 (2.78-2.70)
Ramachandran outliers	100387	3367 (2.78-2.70)
Sidechain outliers	100360	3368 (2.78-2.70)
RSRZ outliers	91569	3055 (2.78-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	454	 4% 60% 32% • 5%
1	B	454	 3% 61% 31% • •

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7209 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 Tyrosine-protein kinase HCK.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	P	S	0	0	0
			3480	2228	583	648	1	20			
1	B	435	Total	C	N	O	P	S	0	0	0
			3507	2242	589	655	1	20			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	GLY	-	EXPRESSION TAG	UNP P08631
A	79	ALA	-	EXPRESSION TAG	UNP P08631
A	80	MET	-	EXPRESSION TAG	UNP P08631
A	81	GLY	-	EXPRESSION TAG	UNP P08631
A	82	SER	-	EXPRESSION TAG	UNP P08631
A	83	GLY	-	EXPRESSION TAG	UNP P08631
A	84	ILE	-	EXPRESSION TAG	UNP P08631
A	85	ARG	-	EXPRESSION TAG	UNP P08631
A	528	GLU	GLN	ENGINEERED MUTATION	UNP P08631
A	529	GLU	GLN	ENGINEERED MUTATION	UNP P08631
A	530	ILE	GLN	ENGINEERED MUTATION	UNP P08631
B	78	GLY	-	EXPRESSION TAG	UNP P08631
B	79	ALA	-	EXPRESSION TAG	UNP P08631
B	80	MET	-	EXPRESSION TAG	UNP P08631
B	81	GLY	-	EXPRESSION TAG	UNP P08631
B	82	SER	-	EXPRESSION TAG	UNP P08631
B	83	GLY	-	EXPRESSION TAG	UNP P08631
B	84	ILE	-	EXPRESSION TAG	UNP P08631
B	85	ARG	-	EXPRESSION TAG	UNP P08631
B	528	GLU	GLN	ENGINEERED MUTATION	UNP P08631
B	529	GLU	GLN	ENGINEERED MUTATION	UNP P08631
B	530	ILE	GLN	ENGINEERED MUTATION	UNP P08631

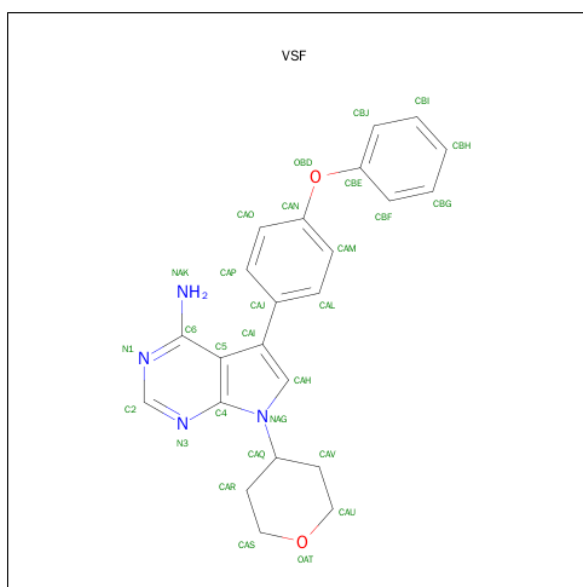
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0
2	A	1	Total Ca 1 1	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	A	1	Total Cl 1 1	0	0

- Molecule 4 is 5-(4-PHENOXYPHENYL)-7-(TETRAHYDRO-2H-PYRAN-4-YL)-7H-PYRR OLO[2,3-D]PYRIMIDIN-4-AMINE (three-letter code: VSF) (formula: C₂₃H₂₂N₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O 29 23 4 2	0	0
4	B	1	Total C N O 29 23 4 2	0	0

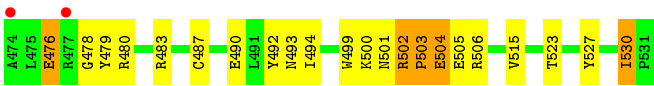
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	84	Total O 84 84	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	76	Total	O	0	0
			76	76		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.81Å 73.75Å 180.13Å 90.00° 95.86° 90.00°	Depositor
Resolution (Å)	40.97 – 2.75 40.97 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.4 (40.97-2.75) 99.4 (40.97-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.205 , 0.265 0.202 , 0.263	Depositor DCC
R_{free} test set	1665 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	48.0	Xtriage
Anisotropy	0.490	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.2	EDS
Estimated twinning fraction	0.026 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 33286 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7209	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CA, PTR, VSF, CL

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	0.24	0/3546	0.42	0/4785
1	B	0.24	0/3574	0.43	0/4825
All	All	0.24	0/7120	0.43	0/9610

There are no bond length outliers.

There are no bond angle outliers.

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	3480	0	3443	154	0
1	B	3507	0	3463	163	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	29	0	22	2	0
4	B	29	0	22	3	0
5	A	84	0	0	7	0
5	B	76	0	0	8	0
All	All	7209	0	6950	315	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 22.

All (315) 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:B:140:ARG:HH11	1:B:140:ARG:HG2	1.12	1.11
1:A:110:MET:HE1	1:A:133:ILE:HG21	1.37	1.04
1:B:110:MET:CE	1:B:133:ILE:HD13	1.93	0.99
1:A:110:MET:HE2	1:A:133:ILE:HD13	1.45	0.98
1:A:221:PHE:HZ	1:A:230:HIS:CD2	1.80	0.98
1:A:191:ASP:OD2	1:A:239:CYS:HB2	1.62	0.97
1:B:140:ARG:HH11	1:B:140:ARG:CG	1.79	0.95
1:B:111:VAL:HG23	1:B:125:LEU:HD21	1.48	0.95
1:A:221:PHE:HZ	1:A:230:HIS:HD2	1.05	0.95
1:B:435:ASN:HB2	1:B:436:PHE:CD2	2.05	0.92
1:B:301:SER:N	1:B:302:MET:HB2	1.88	0.89
1:B:468:ASN:HB2	1:B:469:PRO:HD3	1.54	0.89
1:A:300:GLY:HA3	1:A:302:MET:HE3	1.53	0.88
1:A:191:ASP:OD1	1:A:192:TYR:N	2.07	0.87
1:B:104:PHE:HE2	1:B:133:ILE:HG22	1.40	0.87
1:A:301:SER:N	1:A:302:MET:HB3	1.90	0.87
1:A:426:ILE:HD12	1:A:427:LYS:N	1.89	0.87
1:B:154:ILE:HG22	1:B:176:ARG:HD2	1.57	0.86
1:A:110:MET:CE	1:A:133:ILE:HD13	2.06	0.86
1:B:303:SER:HB2	1:B:306:ALA:HB2	1.58	0.85
1:A:530:ILE:CD1	1:A:531:PRO:HD2	2.06	0.85
1:B:158:ASP:OD1	5:B:716:HOH:O	1.97	0.82
1:A:221:PHE:CZ	1:A:230:HIS:CD2	2.67	0.82
1:B:110:MET:HE3	1:B:133:ILE:HD13	1.61	0.81
1:B:303:SER:HB2	1:B:306:ALA:CB	2.10	0.81
1:A:530:ILE:HD12	1:A:531:PRO:HD2	1.63	0.80
1:A:110:MET:HE1	1:A:133:ILE:CG2	2.09	0.80
1:A:145:GLU:HG3	5:A:739:HOH:O	1.81	0.79
1:B:111:VAL:CG2	1:B:125:LEU:HD21	2.13	0.78
1:A:301:SER:HA	1:A:302:MET:C	2.04	0.78
1:A:178:SER:HB2	1:A:527:PTR:O3P	1.83	0.78
1:B:466:MET:CE	1:B:471:VAL:HA	2.14	0.78
1:B:466:MET:HE2	1:B:471:VAL:HG22	1.66	0.78
1:B:208:LEU:HD21	1:B:212:GLY:HA3	1.67	0.77
1:A:523:THR:O	1:B:490:GLU:HG2	1.85	0.77
1:B:341:MET:CE	1:B:401:LYS:HD2	2.15	0.76
1:B:104:PHE:CE2	1:B:133:ILE:HG22	2.22	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:ARG:NH1	1:B:140:ARG:HG2	1.91	0.75
1:A:232:LYS:O	1:A:241:LYS:NZ	2.20	0.74
1:A:221:PHE:CZ	1:A:230:HIS:HD2	1.97	0.74
1:B:234:GLY:H	1:B:241:LYS:HD2	1.53	0.73
1:A:206:ARG:HH22	1:B:483:ARG:HE	1.34	0.72
1:B:180:THR:HG23	1:B:527:PTR:O1P	1.89	0.72
1:B:341:MET:HE3	1:B:401:LYS:HD2	1.71	0.72
1:A:180:THR:HG21	1:B:493:ASN:ND2	2.06	0.71
1:B:180:THR:OG1	1:B:181:THR:N	2.24	0.70
1:A:175:ILE:HD13	1:A:224:LEU:HD22	1.73	0.70
1:B:435:ASN:HB2	1:B:436:PHE:CE2	2.26	0.70
1:A:90:ALA:HA	1:A:138:VAL:HG12	1.75	0.69
1:A:302:MET:HG2	1:A:302:MET:O	1.93	0.69
1:A:231:TYR:CE1	1:A:236:ASP:HB3	2.28	0.69
1:A:461:ILE:HG13	1:A:462:PRO:HD2	1.74	0.69
1:A:224:LEU:O	1:A:228:VAL:HG23	1.92	0.68
1:A:296:THR:HG22	1:A:335:TYR:CE1	2.28	0.68
1:B:466:MET:HE2	1:B:471:VAL:CG2	2.22	0.68
1:A:380:ARG:NH1	1:A:380:ARG:HB3	2.09	0.68
1:B:466:MET:HE1	1:B:471:VAL:HA	1.76	0.67
1:A:364:ILE:HD12	1:A:520:TYR:OH	1.95	0.67
1:B:302:MET:O	1:B:302:MET:HG2	1.93	0.67
1:A:468:ASN:HB2	1:A:469:PRO:HD3	1.77	0.66
1:A:134:PRO:HG2	1:A:137:TYR:HB2	1.78	0.66
1:A:530:ILE:HD12	1:A:531:PRO:CD	2.26	0.66
1:A:115:GLU:HA	1:A:120:TRP:CD1	2.29	0.66
1:A:193:ASP:OD1	1:A:195:ARG:HG2	1.95	0.66
1:A:380:ARG:HH11	1:A:380:ARG:HB3	1.60	0.66
1:A:360:LEU:HD21	1:A:488:PRO:HD3	1.77	0.66
1:B:397:SER:O	1:B:398:LEU:HB2	1.96	0.66
1:A:181:THR:OG1	1:A:204:LYS:NZ	2.29	0.66
1:B:207:THR:HG23	1:B:213:PHE:CE1	2.30	0.65
1:B:193:ASP:HB3	1:B:196:GLN:H	1.60	0.65
1:A:110:MET:CE	1:A:133:ILE:HG21	2.22	0.65
1:A:300:GLY:HA3	1:A:302:MET:CE	2.24	0.65
1:A:110:MET:HG2	1:A:124:SER:HA	1.79	0.65
1:B:210:ASN:HB2	1:B:211:GLY:HA3	1.78	0.65
1:B:154:ILE:CG2	1:B:176:ARG:HD2	2.25	0.65
1:B:360:LEU:HB3	1:B:361:PRO:HD3	1.79	0.65
1:A:299:PRO:HB3	1:A:332:GLU:HG2	1.80	0.64
1:B:111:VAL:HG23	1:B:125:LEU:CD2	2.25	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:THR:HG21	1:B:493:ASN:HD21	1.63	0.63
1:B:337:ILE:HD12	1:B:337:ILE:N	2.14	0.63
1:A:530:ILE:CG1	1:A:531:PRO:HD2	2.28	0.62
4:B:603:VSF:CAP	4:B:603:VSF:H2	2.11	0.62
1:B:483:ARG:NH1	1:B:487:CYS:O	2.32	0.62
1:A:472:ILE:O	1:A:476:GLU:HG2	1.99	0.62
1:B:237:GLY:O	1:B:530:ILE:HD11	2.00	0.62
1:B:301:SER:CA	1:B:302:MET:HB2	2.30	0.61
1:A:301:SER:H	1:A:302:MET:HB3	1.63	0.61
1:B:503:PRO:O	1:B:506:ARG:N	2.31	0.61
1:B:378:GLU:HG3	1:B:441:ILE:HG12	1.82	0.60
1:A:231:TYR:HE1	1:A:236:ASP:HB3	1.65	0.60
1:A:204:LYS:HD3	1:A:527:PTR:CD2	2.32	0.60
1:A:164:LEU:HA	1:A:190:ARG:NH1	2.17	0.59
1:B:210:ASN:H	1:B:211:GLY:CA	2.15	0.59
1:A:426:ILE:HD11	1:A:468:ASN:OD1	2.02	0.59
1:B:216:SER:HB2	1:B:530:ILE:CG2	2.33	0.59
1:A:530:ILE:HG13	1:A:531:PRO:HD2	1.84	0.59
1:A:461:ILE:HG13	1:A:462:PRO:CD	2.32	0.59
1:B:96:ALA:HA	1:B:101:ASP:OD2	2.02	0.59
1:B:503:PRO:O	1:B:505:GLU:N	2.36	0.59
1:A:490:GLU:HG2	1:B:523:THR:O	2.03	0.58
1:A:476:GLU:OE1	1:A:476:GLU:HA	2.02	0.58
1:B:494:ILE:HD11	1:B:515:VAL:HG11	1.85	0.58
1:A:233:LYS:HB3	5:A:749:HOH:O	2.02	0.58
1:A:227:LEU:C	1:A:227:LEU:HD23	2.24	0.58
1:B:323:VAL:CG1	4:B:603:VSF:H19	2.34	0.58
4:B:603:VSF:H12	5:B:775:HOH:O	2.03	0.57
1:A:403:ALA:O	1:A:404:ASP:HB2	2.03	0.57
1:A:213:PHE:N	1:A:213:PHE:CD1	2.71	0.57
1:A:287:ASN:O	1:A:289:HIS:N	2.37	0.57
1:A:385:ARG:HG3	1:A:439:PHE:CD2	2.41	0.56
1:B:264:ARG:NH2	1:B:331:LYS:O	2.36	0.56
1:A:444:ASP:O	1:A:447:SER:HB2	2.05	0.56
1:B:341:MET:HE1	1:B:401:LYS:HD2	1.85	0.56
1:A:424:PHE:N	1:A:425:PRO:HD3	2.20	0.56
1:B:382:TYR:OH	1:B:406:GLY:HA3	2.05	0.56
1:B:110:MET:HG3	1:B:124:SER:HA	1.88	0.56
1:A:467:SER:OG	1:A:469:PRO:HD2	2.04	0.56
1:B:322:LEU:HD22	1:B:402:ILE:HB	1.88	0.56
1:A:170:LEU:H	1:A:170:LEU:CD1	2.19	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:GLY:CA	1:A:302:MET:HE3	2.32	0.55
1:B:468:ASN:HB2	1:B:469:PRO:CD	2.32	0.55
1:A:432:GLU:OE2	1:A:506:ARG:NH1	2.39	0.55
1:B:140:ARG:NH1	1:B:140:ARG:CG	2.49	0.55
1:A:190:ARG:HG3	1:A:191:ASP:N	2.20	0.55
1:A:269:LEU:N	1:A:269:LEU:HD23	2.21	0.55
1:B:93:ASP:O	1:B:250:LYS:NZ	2.31	0.55
1:B:151:PHE:CD1	1:B:174:MET:HB2	2.41	0.55
1:B:156:ARG:HG3	1:B:202:HIS:CE1	2.42	0.55
1:A:296:THR:HG22	1:A:335:TYR:CD1	2.43	0.54
1:A:163:LEU:HD23	1:A:246:CYS:SG	2.47	0.54
1:B:479:TYR:O	1:B:480:ARG:HD3	2.07	0.54
1:A:189:VAL:HG12	1:A:190:ARG:O	2.07	0.54
1:B:303:SER:HB2	1:B:306:ALA:HB3	1.89	0.54
1:B:208:LEU:HD22	1:B:212:GLY:O	2.08	0.54
1:B:307:PHE:CE2	1:B:334:ILE:HG21	2.43	0.54
1:B:388:ARG:NH1	5:B:774:HOH:O	2.33	0.54
1:B:110:MET:CE	1:B:133:ILE:HG21	2.38	0.54
1:B:301:SER:N	1:B:302:MET:CB	2.66	0.54
1:B:387:LEU:O	1:B:388:ARG:HB3	2.08	0.54
1:B:301:SER:HB2	1:B:302:MET:HB2	1.89	0.53
1:A:337:ILE:HD12	1:A:337:ILE:N	2.23	0.53
1:A:221:PHE:CD1	1:A:227:LEU:HA	2.43	0.53
1:B:154:ILE:CG2	1:B:176:ARG:CD	2.86	0.53
1:B:210:ASN:H	1:B:211:GLY:HA3	1.72	0.53
1:A:442:LYS:NZ	1:A:506:ARG:O	2.36	0.53
1:B:426:ILE:HD12	1:B:426:ILE:C	2.29	0.53
1:B:530:ILE:O	1:B:530:ILE:HG22	2.08	0.53
1:A:503:PRO:HD2	1:A:504:GLU:OE2	2.09	0.53
1:A:467:SER:O	1:A:471:VAL:HG23	2.09	0.53
1:B:397:SER:O	1:B:398:LEU:CB	2.57	0.53
1:A:164:LEU:HD21	1:A:200:VAL:HG23	1.91	0.52
1:B:501:ASN:O	1:B:502:ARG:CB	2.57	0.52
1:B:104:PHE:CE2	1:B:133:ILE:CG2	2.92	0.52
1:B:222:SER:HB3	1:B:226:GLU:OE2	2.09	0.52
1:B:441:ILE:O	1:B:445:VAL:HG23	2.09	0.52
1:A:304:VAL:CG2	1:A:305:GLU:N	2.73	0.52
1:A:377:ILE:HG23	1:A:382:TYR:HB3	1.91	0.52
1:A:433:ALA:HA	1:A:438:SER:H	1.75	0.52
1:A:301:SER:CA	1:A:302:MET:HB3	2.40	0.52
1:B:436:PHE:N	1:B:436:PHE:CD2	2.77	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:504:GLU:H	1:A:504:GLU:CD	2.14	0.51
1:A:486:ASN:N	1:A:486:ASN:OD1	2.41	0.51
1:B:424:PHE:CD2	1:B:424:PHE:N	2.78	0.51
1:B:150:PHE:HE2	1:B:152:LYS:HD2	1.76	0.51
1:A:158:ASP:HB2	5:A:721:HOH:O	2.10	0.51
1:A:156:ARG:HG3	1:A:202:HIS:CE1	2.45	0.51
1:B:172:SER:HA	1:B:244:VAL:O	2.10	0.51
1:B:208:LEU:HD22	1:B:212:GLY:C	2.31	0.51
1:B:308:LEU:CD2	1:B:330:THR:HG22	2.40	0.51
1:A:170:LEU:HD12	1:A:170:LEU:N	2.26	0.51
1:A:442:LYS:HE2	5:A:701:HOH:O	2.11	0.51
1:A:411:ILE:HG23	1:A:412:GLU:N	2.26	0.51
1:B:426:ILE:HD12	1:B:427:LYS:N	2.26	0.51
1:B:210:ASN:CB	1:B:211:GLY:HA3	2.35	0.51
1:B:504:GLU:CD	1:B:504:GLU:H	2.14	0.51
1:A:404:ASP:OD1	4:A:603:VSF:H15	2.10	0.51
1:B:277:GLN:O	1:B:298:LYS:HE2	2.11	0.51
1:B:210:ASN:N	1:B:211:GLY:HA3	2.25	0.50
1:B:502:ARG:HG2	1:B:504:GLU:OE2	2.10	0.50
1:B:86:ILE:C	1:B:87:ILE:HD12	2.32	0.50
1:B:430:ALA:HB2	1:B:446:TRP:CB	2.42	0.50
1:A:389:ALA:N	1:A:451:LEU:HD13	2.27	0.50
1:B:453:MET:HA	1:B:456:VAL:HG12	1.94	0.50
1:B:87:ILE:HD12	1:B:87:ILE:N	2.27	0.49
1:A:303:SER:HB3	1:A:306:ALA:CB	2.42	0.49
1:A:109:GLN:C	1:A:110:MET:HG3	2.33	0.49
1:B:137:TYR:HE2	1:B:250:LYS:HE3	1.78	0.49
1:A:452:LEU:HD23	1:A:495:MET:HG2	1.93	0.49
1:B:468:ASN:CB	1:B:469:PRO:HD3	2.33	0.49
1:B:308:LEU:HD21	1:B:330:THR:HG22	1.94	0.49
1:B:263:PRO:O	1:B:266:SER:OG	2.28	0.49
1:A:277:GLN:O	1:A:298:LYS:HE2	2.13	0.49
1:B:87:ILE:CD1	1:B:87:ILE:N	2.76	0.48
1:B:208:LEU:CD2	1:B:212:GLY:HA3	2.39	0.48
1:B:154:ILE:HG21	1:B:176:ARG:HD3	1.94	0.48
1:B:504:GLU:CD	1:B:504:GLU:N	2.67	0.48
1:A:230:HIS:ND1	1:A:230:HIS:C	2.67	0.48
1:B:115:GLU:HA	1:B:120:TRP:CD1	2.48	0.48
1:B:302:MET:O	1:B:303:SER:C	2.52	0.48
1:A:170:LEU:N	1:A:170:LEU:CD1	2.76	0.48
1:B:424:PHE:N	1:B:425:PRO:HD3	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LEU:HG	1:A:131:GLY:HA3	1.94	0.47
1:A:529:GLU:HG3	5:B:717:HOH:O	2.14	0.47
1:B:208:LEU:HD22	1:B:208:LEU:H	1.79	0.47
1:B:110:MET:HE1	1:B:133:ILE:HG21	1.96	0.47
1:B:154:ILE:HG23	1:B:155:SER:O	2.14	0.47
1:A:141:VAL:C	1:A:143:SER:H	2.18	0.47
1:B:303:SER:O	1:B:306:ALA:HB3	2.13	0.47
1:B:347:LEU:O	1:B:351:LYS:HG3	2.13	0.47
1:B:86:ILE:HD13	1:B:140:ARG:NH2	2.29	0.47
1:B:343:LYS:HE3	1:B:396:ALA:O	2.14	0.47
1:B:424:PHE:N	1:B:425:PRO:CD	2.77	0.47
1:B:446:TRP:CE3	1:B:499:TRP:HA	2.50	0.47
1:B:253:LYS:HD3	5:B:736:HOH:O	2.13	0.47
1:B:210:ASN:N	1:B:211:GLY:CA	2.77	0.47
1:A:303:SER:HB3	1:A:306:ALA:HB2	1.95	0.47
1:A:134:PRO:HG2	1:A:137:TYR:CB	2.45	0.47
1:A:302:MET:CG	1:A:302:MET:O	2.61	0.47
1:B:208:LEU:CD2	1:B:212:GLY:CA	2.92	0.46
1:B:265:GLU:CD	1:B:265:GLU:H	2.18	0.46
1:A:215:ILE:HG22	1:A:216:SER:N	2.30	0.46
1:B:222:SER:HB3	1:B:226:GLU:CD	2.36	0.46
1:A:104:PHE:HE2	1:A:133:ILE:HG22	1.81	0.46
1:A:384:HIS:O	1:A:385:ARG:HB3	2.16	0.46
1:B:301:SER:CB	1:B:302:MET:HB2	2.46	0.46
1:B:483:ARG:HB2	1:B:492:TYR:CD1	2.50	0.45
1:B:476:GLU:C	1:B:478:GLY:H	2.20	0.45
1:A:204:LYS:HD3	1:A:527:PTR:CG	2.44	0.45
1:A:291:LYS:HE2	1:A:340:PHE:CD2	2.52	0.45
1:B:303:SER:O	1:B:306:ALA:N	2.49	0.45
1:A:173:PHE:O	1:A:174:MET:HB3	2.15	0.45
4:A:603:VSF:H2	4:A:603:VSF:CAP	2.29	0.45
1:B:467:SER:OG	1:B:469:PRO:HD2	2.17	0.45
1:A:408:ALA:HB2	5:A:717:HOH:O	2.17	0.45
1:B:304:VAL:HG21	1:B:332:GLU:HG3	1.99	0.45
1:B:206:ARG:HB2	1:B:214:TYR:CE1	2.52	0.45
1:B:503:PRO:HA	1:B:506:ARG:HD2	1.97	0.45
1:A:433:ALA:O	1:A:437:GLY:HA2	2.16	0.45
1:A:98:HIS:HE1	5:A:748:HOH:O	2.00	0.44
1:B:385:ARG:HG3	1:B:439:PHE:CD2	2.53	0.44
1:A:296:THR:CG2	1:A:335:TYR:CE1	3.00	0.44
1:B:430:ALA:HB2	1:B:446:TRP:HB3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:300:GLY:C	1:B:302:MET:CB	2.86	0.44
1:A:140:ARG:O	1:A:143:SER:HB3	2.18	0.44
1:A:94:TYR:HB3	1:A:104:PHE:CE1	2.52	0.44
1:B:210:ASN:H	1:B:211:GLY:C	2.21	0.44
1:B:115:GLU:HG2	1:B:120:TRP:CE2	2.52	0.44
1:B:102:LEU:HG	1:B:131:GLY:HA3	1.98	0.44
1:B:304:VAL:O	1:B:308:LEU:HG	2.18	0.44
1:B:388:ARG:HD2	5:B:774:HOH:O	2.17	0.44
1:A:104:PHE:CE2	1:A:133:ILE:HG22	2.53	0.44
1:A:220:THR:O	1:A:221:PHE:CD2	2.70	0.44
1:B:303:SER:CB	1:B:306:ALA:HB2	2.38	0.44
1:A:302:MET:SD	5:A:743:HOH:O	2.61	0.43
1:A:178:SER:OG	1:A:181:THR:O	2.34	0.43
1:A:364:ILE:HD13	1:A:519:PHE:CD2	2.53	0.43
1:B:269:LEU:HD23	1:B:269:LEU:N	2.33	0.43
1:A:446:TRP:CE3	1:A:499:TRP:HA	2.54	0.43
1:A:190:ARG:HG3	1:A:191:ASP:H	1.82	0.43
1:B:451:LEU:HD12	1:B:451:LEU:O	2.19	0.43
1:B:424:PHE:HD2	1:B:424:PHE:N	2.17	0.43
1:A:530:ILE:HG13	1:A:531:PRO:CD	2.48	0.43
1:B:161:ARG:NH1	5:B:716:HOH:O	2.52	0.43
1:A:127:THR:O	1:A:128:ARG:CB	2.67	0.43
1:A:360:LEU:HD21	1:A:488:PRO:CD	2.47	0.43
1:A:109:GLN:O	1:A:110:MET:HG3	2.19	0.43
1:A:426:ILE:C	1:A:426:ILE:HD12	2.37	0.43
1:B:383:ILE:HG22	1:B:384:HIS:N	2.33	0.43
1:B:466:MET:HE2	1:B:471:VAL:HA	1.96	0.42
1:A:476:GLU:CA	1:A:476:GLU:OE1	2.67	0.42
1:A:300:GLY:O	1:A:301:SER:CB	2.66	0.42
1:A:468:ASN:N	1:A:469:PRO:CD	2.82	0.42
1:A:325:LEU:HD12	1:A:325:LEU:N	2.34	0.42
1:A:134:PRO:HG2	1:A:137:TYR:CG	2.54	0.42
1:B:387:LEU:HA	1:B:387:LEU:HD23	1.84	0.42
1:B:301:SER:CA	1:B:302:MET:CB	2.95	0.42
1:A:204:LYS:C	1:A:205:ILE:HG13	2.40	0.42
1:B:440:THR:N	1:B:443:SER:OG	2.50	0.42
1:A:479:TYR:O	1:A:480:ARG:HD3	2.20	0.42
1:A:341:MET:SD	1:A:401:LYS:HD2	2.60	0.42
1:A:88:VAL:HG13	1:A:112:VAL:HG23	2.02	0.42
1:A:160:GLU:O	1:A:164:LEU:HG	2.20	0.41
1:B:260:TRP:CZ3	1:B:315:LYS:HD3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:426:ILE:HD12	1:A:427:LYS:H	1.74	0.41
1:B:432:GLU:CD	1:B:432:GLU:H	2.24	0.41
1:B:377:ILE:HG23	1:B:382:TYR:HB3	2.02	0.41
1:A:502:ARG:HB3	1:A:504:GLU:OE2	2.19	0.41
1:B:279:GLY:HA3	1:B:296:THR:O	2.20	0.41
1:A:257:LYS:HG3	1:A:257:LYS:H	1.65	0.41
1:B:476:GLU:C	1:B:478:GLY:N	2.73	0.41
1:B:349:PHE:O	1:B:355:GLY:HA3	2.21	0.41
1:B:301:SER:CB	1:B:302:MET:CA	2.98	0.41
1:A:204:LYS:O	1:A:205:ILE:HG13	2.20	0.41
1:A:502:ARG:HA	1:A:503:PRO:HD3	1.94	0.41
1:B:223:THR:OG1	1:B:226:GLU:HG3	2.21	0.41
1:A:287:ASN:C	1:A:289:HIS:N	2.74	0.41
1:B:427:LYS:HD3	1:B:463:TYR:HD1	1.86	0.41
1:A:222:SER:HB3	1:A:226:GLU:OE2	2.20	0.41
1:B:255:TRP:CD2	1:B:326:HIS:HB3	2.56	0.41
1:A:387:LEU:HD23	1:A:387:LEU:HA	1.85	0.41
1:A:426:ILE:HG12	1:A:468:ASN:HB3	2.02	0.41
1:A:287:ASN:C	1:A:289:HIS:H	2.24	0.41
1:A:432:GLU:H	1:A:432:GLU:HG3	1.46	0.41
1:B:300:GLY:C	1:B:302:MET:HB3	2.41	0.40
1:A:530:ILE:CG1	1:A:531:PRO:CD	2.99	0.40
1:A:102:LEU:HD11	1:A:131:GLY:N	2.36	0.40
1:A:493:ASN:O	1:A:497:ARG:HG3	2.21	0.40
1:B:500:LYS:HA	1:B:500:LYS:HD3	1.87	0.40
1:A:264:ARG:NH2	1:A:331:LYS:O	2.51	0.40
1:A:426:ILE:HD11	1:A:468:ASN:CG	2.40	0.40
1:B:154:ILE:HG21	1:B:176:ARG:CD	2.51	0.40
1:B:98:HIS:HB3	5:B:728:HOH:O	2.21	0.40
1:A:97:ILE:HG13	1:A:98:HIS:CE1	2.56	0.40
1:B:503:PRO:HB2	1:B:504:GLU:H	1.77	0.40
1:B:262:ILE:HB	1:B:263:PRO:HD2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	423/454 (93%)	388 (92%)	32 (8%)	3 (1%)	26	54
1	B	430/454 (95%)	388 (90%)	32 (7%)	10 (2%)	8	18
All	All	853/908 (94%)	776 (91%)	64 (8%)	13 (2%)	13	30

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	301	SER
1	B	302	MET
1	B	410	VAL
1	B	504	GLU
1	A	234	GLY
1	A	288	LYS
1	B	234	GLY
1	B	301	SER
1	B	303	SER
1	B	356	SER
1	B	398	LEU
1	B	503	PRO
1	B	502	ARG

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	376/393 (96%)	354 (94%)	22 (6%)	24	49
1	B	378/393 (96%)	356 (94%)	22 (6%)	25	50
All	All	754/786 (96%)	710 (94%)	44 (6%)	25	50

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	141	VAL
1	A	178	SER
1	A	181	THR
1	A	204	LYS
1	A	213	PHE
1	A	240	GLN
1	A	244	VAL
1	A	249	SER
1	A	257	LYS
1	A	280	GLU
1	A	297	MET
1	A	302	MET
1	A	304	VAL
1	A	328	VAL
1	A	387	LEU
1	A	432	GLU
1	A	456	VAL
1	A	461	ILE
1	A	501	ASN
1	A	504	GLU
1	A	505	GLU
1	A	514	SER
1	B	92	TYR
1	B	110	MET
1	B	140	ARG
1	B	180	THR
1	B	184	SER
1	B	193	ASP
1	B	195	ARG
1	B	208	LEU
1	B	209	ASP
1	B	218	ARG
1	B	220	THR
1	B	249	SER
1	B	253	LYS
1	B	305	GLU
1	B	352	SER
1	B	353	ASP
1	B	387	LEU
1	B	424	PHE
1	B	432	GLU
1	B	467	SER
1	B	476	GLU

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Mol	Chain	Res	Type
1	B	530	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	230	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	PTR	A	527	1,2	14,16,17	1.31	1 (7%)	18,22,24	0.71	0
1	PTR	B	527	1,2	14,16,17	1.26	1 (7%)	18,22,24	0.67	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
1	PTR	A	527	1,2	-	0/9/11/13	0/1/1/1
1	PTR	B	527	1,2	-	0/9/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	527	PTR	OH-CZ	-4.34	1.30	1.40
1	B	527	PTR	OH-CZ	-4.09	1.30	1.40

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	527	PTR	3	0
1	B	527	PTR	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
4	VSF	A	603	-	31,33,33	0.73	1 (3%)	35,46,46	1.94	2 (5%)
4	VSF	B	603	-	31,33,33	0.76	1 (3%)	35,46,46	1.97	2 (5%)

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
4	VSF	A	603	-	-	0/8/20/20	0/5/5/5
4	VSF	B	603	-	-	0/8/20/20	0/5/5/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	603	VSF	CAH-NAG	-2.49	1.34	1.38
4	A	603	VSF	CAH-NAG	-2.44	1.34	1.38

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	603	VSF	N3-C2-N1	-10.61	120.77	128.89
4	A	603	VSF	N3-C2-N1	-10.42	120.92	128.89
4	A	603	VSF	CBE-OBD-CAN	2.03	123.87	118.81
4	B	603	VSF	CBE-OBD-CAN	2.58	125.25	118.81

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	603	VSF	2	0
4	B	603	VSF	3	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	429/454 (94%)	0.15	16 (3%) 45 47	25, 44, 85, 107	0
1	B	434/454 (95%)	0.13	13 (2%) 54 56	28, 47, 82, 103	0
All	All	863/908 (95%)	0.14	29 (3%) 49 51	25, 46, 84, 107	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	VAL	4.5
1	B	303	SER	4.2
1	B	305	GLU	3.9
1	B	474	ALA	3.9
1	A	301	SER	3.9
1	B	211	GLY	3.9
1	A	305	GLU	3.7
1	B	301	SER	3.6
1	A	234	GLY	3.4
1	B	469	PRO	3.3
1	A	142	ASP	3.2
1	A	216	SER	3.2
1	B	436	PHE	2.9
1	A	219	SER	2.9
1	A	531	PRO	2.7
1	A	233	LYS	2.7
1	A	214	TYR	2.6
1	A	465	GLY	2.5
1	A	223	THR	2.5
1	B	473	ARG	2.4
1	A	237	GLY	2.4
1	B	465	GLY	2.3
1	A	224	LEU	2.3
1	B	432	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	228	VAL	2.1
1	B	477	ARG	2.1
1	B	300	GLY	2.1
1	A	213	PHE	2.0
1	A	215	ILE	2.0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
1	PTR	A	527	16/17	0.97	0.16	-	38,49,55,56	0
1	PTR	B	527	16/17	0.97	0.18	-	24,40,44,45	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	VSF	A	603	29/29	0.97	0.19	0.71	22,28,34,44	0
4	VSF	B	603	29/29	0.98	0.18	0.30	26,32,41,47	0
2	CA	B	601	1/1	0.96	0.17	-0.22	48,48,48,48	0
3	CL	B	602	1/1	0.99	0.16	-	47,47,47,47	0
3	CL	A	602	1/1	0.99	0.13	-	55,55,55,55	0
2	CA	A	601	1/1	0.90	0.22	-	49,49,49,49	0

6.5 Other polymers

There are no such residues in this entry.