



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:06 PM UTC

PDB ID : 8BE6 / pdb_00008be6
Title : Crystal structure of SOS1-HRas-peptidomimetic2
Authors : Fischer, B.; Wohlkonig, A.; Steyaert, J.
Deposited on : 2022-10-21
Resolution : 2.90 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)	:	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

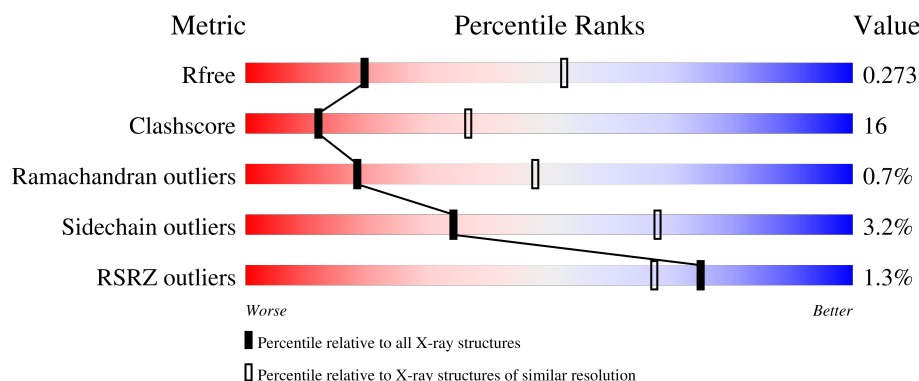
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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 $\leq 5\%$. 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	R	186	% 68% 19% • 11%
2	S	507	% 59% 26% • 13%
3	P	8	12% 12% 50% 25%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5129 atoms, of which 58 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 GTPase HRas.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	R	166	Total	C	N	O	S	3	0	0
			1322	824	228	263	7			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	-19	MET	-	initiating methionine	UNP P01112
R	-18	GLY	-	expression tag	UNP P01112
R	-17	SER	-	expression tag	UNP P01112
R	-16	SER	-	expression tag	UNP P01112
R	-15	HIS	-	expression tag	UNP P01112
R	-14	HIS	-	expression tag	UNP P01112
R	-13	HIS	-	expression tag	UNP P01112
R	-12	HIS	-	expression tag	UNP P01112
R	-11	HIS	-	expression tag	UNP P01112
R	-10	HIS	-	expression tag	UNP P01112
R	-9	SER	-	expression tag	UNP P01112
R	-8	SER	-	expression tag	UNP P01112
R	-7	GLY	-	expression tag	UNP P01112
R	-6	LEU	-	expression tag	UNP P01112
R	-5	VAL	-	expression tag	UNP P01112
R	-4	PRO	-	expression tag	UNP P01112
R	-3	ARG	-	expression tag	UNP P01112
R	-2	GLY	-	expression tag	UNP P01112
R	-1	SER	-	expression tag	UNP P01112
R	0	HIS	-	expression tag	UNP P01112

- Molecule 2 is a protein called Son of sevenless homolog 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	S	440	Total	C	N	O	S	79	0	0
			3668	2364	629	663	12			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	543	MET	-	initiating methionine	UNP Q07889
S	544	GLY	-	expression tag	UNP Q07889
S	545	SER	-	expression tag	UNP Q07889
S	546	SER	-	expression tag	UNP Q07889
S	547	HIS	-	expression tag	UNP Q07889
S	548	HIS	-	expression tag	UNP Q07889
S	549	HIS	-	expression tag	UNP Q07889
S	550	HIS	-	expression tag	UNP Q07889
S	551	HIS	-	expression tag	UNP Q07889
S	552	HIS	-	expression tag	UNP Q07889
S	553	SER	-	expression tag	UNP Q07889
S	554	SER	-	expression tag	UNP Q07889
S	555	GLY	-	expression tag	UNP Q07889
S	556	LEU	-	expression tag	UNP Q07889
S	557	VAL	-	expression tag	UNP Q07889
S	558	PRO	-	expression tag	UNP Q07889
S	559	ARG	-	expression tag	UNP Q07889
S	560	GLY	-	expression tag	UNP Q07889
S	561	SER	-	expression tag	UNP Q07889
S	562	HIS	-	expression tag	UNP Q07889
S	563	MET	-	expression tag	UNP Q07889

- Molecule 3 is a protein called SOS1-HRas-peptidomimetic2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	8	Total	C	H	N	O	0	0	1
			119	39	58	14	8			

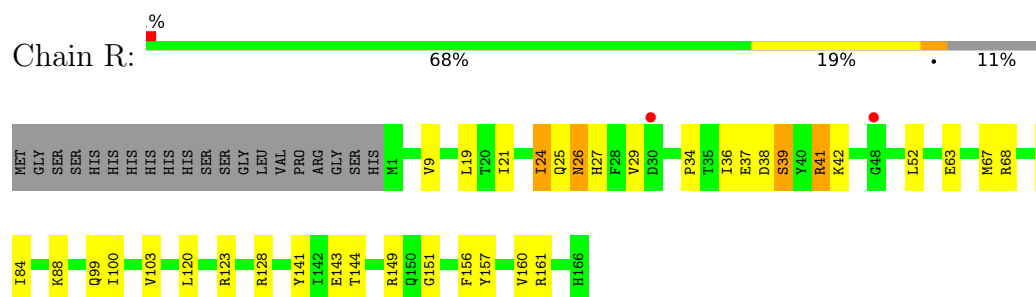
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	R	11	Total	O	0	0
			11	11		
4	S	9	Total	O	0	0
			9	9		

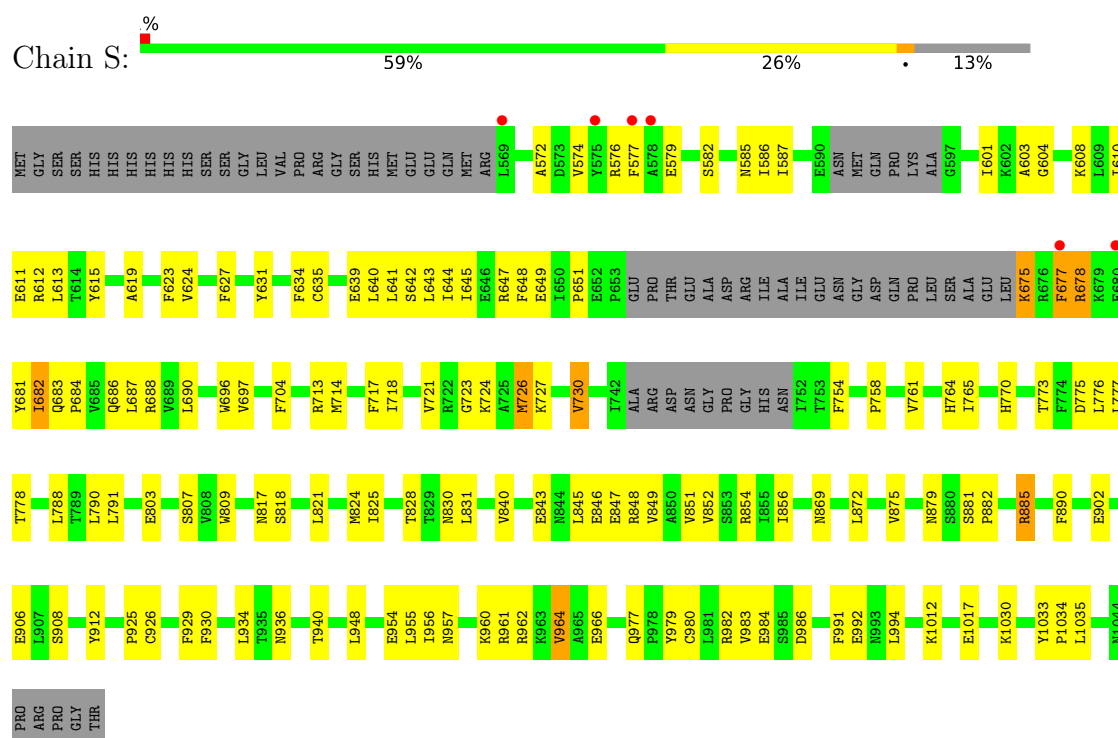
3 Residue-property plots [i](#)

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: GTPase HRas



- Molecule 2: Son of sevenless homolog 1



- Molecule 3: SOS1-HRas-peptidomimetic2



R1	R2	P3	W4	S5	V6	A7	RH28
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4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	142.32Å 142.32Å 207.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	72.25 – 2.90 72.25 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (72.25-2.90) 99.0 (72.25-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.23 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.220 , 0.273 (Not available) , 0.273	Depositor DCC
R_{free} test set	1190 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	94.8	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.010 for -1/2*h-1/2*k-1/2*l,-1/2*h-1/2*k+1/2*l,-h+k 0.003 for -1/2*h+1/2*k-1/2*l,1/2*h-1/2*k-1/2*l,-h-k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5129	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.04% 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: DAR, NH2

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	R	0.23	0/1341	0.44	1/1810 (0.1%)
2	S	0.17	0/3757	0.37	0/5080
3	P	5.07	10/52 (19.2%)	1.96	3/72 (4.2%)
All	All	0.54	10/5150 (0.2%)	0.44	4/6962 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	3	PRO	C-N	14.14	1.51	1.33
3	P	4	TRP	C-N	13.30	1.52	1.33
3	P	3	PRO	CA-CB	-12.86	1.35	1.53
3	P	5	SER	C-N	12.13	1.50	1.33
3	P	3	PRO	N-CA	12.11	1.62	1.47
3	P	6	VAL	C-N	11.15	1.49	1.33
3	P	3	PRO	N-CD	-9.97	1.33	1.47
3	P	2	HIS	C-N	7.50	1.51	1.33
3	P	2	HIS	CB-CG	5.33	1.57	1.50
3	P	2	HIS	CD2-NE2	-5.00	1.32	1.37

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	6	VAL	O-C-N	-5.47	115.73	122.57
1	R	73	ARG	CB-CG-CD	-5.28	99.15	111.30
3	P	2	HIS	CA-C-N	5.01	126.10	119.84
3	P	2	HIS	C-N-CA	5.01	126.10	119.84

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	R	1322	0	1296	33	0
2	S	3668	0	3699	122	0
3	P	61	58	57	4	0
4	R	11	0	0	1	0
4	S	9	0	0	0	0
All	All	5071	58	5052	155	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 16.

All (155) 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)
2:S:610:ILE:HD13	2:S:643:LEU:HD13	1.45	0.96
2:S:678:ARG:HH11	2:S:678:ARG:HB3	1.36	0.91
2:S:936:ASN:O	2:S:940:THR:HG23	1.75	0.87
2:S:681:TYR:HD1	2:S:682:ILE:HD12	1.46	0.79
2:S:840:VAL:CG1	2:S:1012:LYS:HB3	2.19	0.73
1:R:9:VAL:HG21	1:R:100:ILE:HD11	1.71	0.72
2:S:610:ILE:HG21	2:S:643:LEU:HD13	1.71	0.71
2:S:675:LYS:HB2	2:S:675:LYS:NZ	2.06	0.70
2:S:613:LEU:CD1	2:S:624:VAL:HG22	2.22	0.69
2:S:613:LEU:HD11	2:S:624:VAL:HG22	1.74	0.69
2:S:686:GLN:O	2:S:690:LEU:HD12	1.90	0.69
1:R:36:ILE:HG22	1:R:38:ASP:H	1.56	0.69
2:S:925:PRO:HD3	2:S:977:GLN:HG2	1.74	0.69
2:S:610:ILE:HG21	2:S:643:LEU:CD1	2.23	0.69
2:S:791:LEU:HD11	2:S:992:GLU:HG3	1.77	0.66
2:S:721:VAL:HG21	2:S:730:VAL:HG21	1.79	0.65
2:S:714:MET:O	2:S:718:ILE:HG13	1.98	0.64
2:S:678:ARG:HH11	2:S:678:ARG:CB	2.10	0.64
2:S:681:TYR:CD1	2:S:682:ILE:HD12	2.33	0.63
1:R:149:ARG:NH1	4:R:201:HOH:O	2.28	0.62
2:S:684:PRO:O	2:S:688:ARG:HG2	1.99	0.62
2:S:847:GLU:HG2	2:S:1035:LEU:HD21	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:803:GLU:HG2	2:S:818:SER:HA	1.82	0.61
1:R:21:ILE:O	1:R:25:GLN:HB3	2.02	0.59
2:S:777:LEU:HA	2:S:854:ARG:NH1	2.16	0.59
2:S:777:LEU:CD2	2:S:854:ARG:HG3	2.33	0.59
2:S:579:GLU:O	2:S:608:LYS:NZ	2.33	0.59
2:S:961:ARG:O	2:S:964:VAL:HG22	2.02	0.59
2:S:718:ILE:HG22	2:S:730:VAL:CG1	2.33	0.59
2:S:846:GLU:O	2:S:849:VAL:HG22	2.03	0.59
2:S:885:ARG:NH2	2:S:1017:GLU:O	2.36	0.58
1:R:26:ASN:O	1:R:26:ASN:OD1	2.23	0.57
2:S:960:LYS:O	2:S:964:VAL:HG13	2.04	0.57
2:S:576:ARG:NH2	2:S:647:ARG:HH11	2.03	0.57
2:S:645:ILE:O	2:S:649:GLU:HG3	2.05	0.57
2:S:635:CYS:SG	2:S:639:GLU:HG3	2.45	0.56
2:S:788:LEU:HD21	2:S:991:PHE:CD1	2.40	0.56
1:R:84:ILE:HG23	1:R:123:ARG:HG3	1.87	0.56
2:S:770:HIS:HB3	2:S:773:THR:HG22	1.86	0.56
2:S:718:ILE:HG22	2:S:730:VAL:HG11	1.88	0.56
2:S:790:LEU:HD13	2:S:983:VAL:CG2	2.36	0.56
2:S:675:LYS:HD3	2:S:677:PHE:HB2	1.87	0.55
2:S:984:GLU:OE1	2:S:986:ASP:N	2.39	0.55
2:S:641:LEU:O	2:S:644:ILE:HG13	2.07	0.54
2:S:683:GLN:O	2:S:687:LEU:HD23	2.08	0.54
2:S:718:ILE:CG2	2:S:730:VAL:HG13	2.38	0.54
1:R:26:ASN:O	1:R:29:VAL:HB	2.08	0.53
2:S:585:ASN:HB3	2:S:608:LYS:HD3	1.89	0.53
2:S:872:LEU:HA	2:S:875:VAL:HG22	1.90	0.53
2:S:577:PHE:HZ	2:S:647:ARG:HB2	1.72	0.53
2:S:648:PHE:CE1	2:S:686:GLN:HB3	2.43	0.53
2:S:640:LEU:HD23	2:S:696:TRP:CH2	2.44	0.53
2:S:845:LEU:HD23	2:S:845:LEU:O	2.09	0.53
1:R:26:ASN:OD1	1:R:26:ASN:C	2.52	0.52
2:S:634:PHE:CD1	2:S:956:ILE:HD13	2.44	0.52
2:S:824:MET:O	2:S:828:THR:HG23	2.08	0.52
1:R:157:TYR:O	1:R:161:ARG:HG3	2.09	0.52
2:S:803:GLU:HG2	2:S:817:ASN:O	2.10	0.52
2:S:845:LEU:HD23	2:S:845:LEU:C	2.35	0.52
2:S:601:ILE:HD11	2:S:612:ARG:HG3	1.90	0.52
2:S:640:LEU:HD12	2:S:643:LEU:HD11	1.92	0.52
2:S:761:VAL:HG23	2:S:982:ARG:HD3	1.92	0.51
2:S:790:LEU:HD13	2:S:983:VAL:HG23	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:603:ALA:HB2	2:S:948:LEU:HD12	1.92	0.51
2:S:957:ASN:O	2:S:961:ARG:NH1	2.36	0.51
2:S:758:PRO:O	2:S:982:ARG:NH1	2.43	0.51
2:S:645:ILE:HG12	2:S:717:PHE:CD1	2.46	0.51
1:R:34:PRO:HA	1:R:39:SER:OG	2.11	0.51
2:S:686:GLN:OE1	2:S:726:MET:HE1	2.10	0.51
1:R:27:HIS:C	1:R:29:VAL:H	2.20	0.50
2:S:613:LEU:HB2	2:S:623:PHE:HE2	1.77	0.50
2:S:830:ASN:ND2	2:S:994:LEU:O	2.44	0.50
2:S:890:PHE:CG	3:P:3:PRO:HB2	2.47	0.50
1:R:36:ILE:HD12	1:R:36:ILE:N	2.26	0.50
2:S:902:GLU:O	2:S:906:GLU:HG3	2.11	0.50
2:S:576:ARG:HH22	2:S:647:ARG:HD2	1.77	0.50
2:S:809:TRP:CZ2	2:S:934:LEU:HB3	2.46	0.50
1:R:63:GLU:HB3	2:S:825:ILE:HG21	1.93	0.49
1:R:24:ILE:HA	1:R:42:LYS:HE3	1.94	0.49
2:S:776:LEU:O	2:S:776:LEU:HD23	2.12	0.49
1:R:156:PHE:O	1:R:160:VAL:HG23	2.12	0.49
2:S:577:PHE:CE1	2:S:647:ARG:HD3	2.48	0.49
2:S:954:GLU:C	2:S:955:LEU:HD23	2.38	0.49
2:S:619:ALA:HB2	2:S:688:ARG:CZ	2.43	0.49
2:S:908:SER:HA	2:S:912:TYR:CD1	2.48	0.48
2:S:697:VAL:CG1	2:S:714:MET:HE2	2.43	0.48
2:S:847:GLU:O	2:S:851:VAL:HG23	2.13	0.48
2:S:586:ILE:HD11	2:S:608:LYS:HB3	1.94	0.48
2:S:675:LYS:HB2	2:S:675:LYS:HZ2	1.78	0.48
2:S:608:LYS:NZ	2:S:611:GLU:OE1	2.40	0.48
2:S:714:MET:CE	2:S:718:ILE:HD11	2.43	0.48
2:S:641:LEU:HD12	2:S:644:ILE:HD11	1.95	0.48
1:R:141:TYR:CZ	1:R:143:GLU:HG2	2.49	0.48
2:S:962:ARG:O	2:S:966:GLU:HG3	2.14	0.48
2:S:872:LEU:HD12	2:S:929:PHE:CG	2.49	0.48
2:S:601:ILE:O	2:S:962:ARG:NH2	2.26	0.48
1:R:67:MET:HE3	1:R:71:TYR:CE2	2.49	0.47
2:S:764:HIS:CE1	2:S:765:ILE:HG13	2.49	0.47
1:R:72:MET:HE3	1:R:100:ILE:HA	1.95	0.47
2:S:840:VAL:HG13	2:S:1012:LYS:HB3	1.97	0.46
2:S:881:SER:HB2	2:S:882:PRO:HD3	1.97	0.46
2:S:714:MET:HE1	2:S:718:ILE:HD11	1.97	0.46
3:P:6:VAL:O	3:P:6:VAL:HG22	2.15	0.46
2:S:847:GLU:CD	2:S:1030:LYS:H	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:613:LEU:O	2:S:613:LEU:HD12	2.16	0.46
1:R:41:ARG:HH21	1:R:52:LEU:HD21	1.81	0.46
2:S:724:LYS:HA	2:S:727:LYS:HE2	1.97	0.46
1:R:72:MET:HE1	1:R:99:GLN:HG2	1.98	0.46
2:S:872:LEU:HD12	2:S:929:PHE:CD1	2.50	0.46
3:P:2:HIS:HB3	3:P:4:TRP:CE2	2.51	0.46
2:S:704:PHE:CE1	2:S:714:MET:HG3	2.51	0.45
2:S:807:SER:HA	2:S:809:TRP:CZ3	2.50	0.45
2:S:587:ILE:HD11	2:S:955:LEU:HD12	1.99	0.45
2:S:764:HIS:ND1	2:S:765:ILE:HG13	2.32	0.45
2:S:681:TYR:O	2:S:684:PRO:HD2	2.17	0.44
2:S:723:GLY:O	2:S:727:LYS:HG3	2.16	0.44
1:R:24:ILE:HD11	1:R:39:SER:O	2.16	0.44
2:S:926:CYS:HA	2:S:979:TYR:OH	2.17	0.44
2:S:642:SER:HA	2:S:713:ARG:HH22	1.82	0.44
1:R:99:GLN:O	1:R:103:VAL:HG13	2.18	0.44
2:S:582:SER:H	2:S:585:ASN:HB2	1.83	0.44
2:S:775:ASP:N	2:S:778:THR:OG1	2.44	0.44
2:S:925:PRO:CD	2:S:977:GLN:HG2	2.46	0.44
2:S:608:LYS:HA	2:S:611:GLU:HB3	2.00	0.43
2:S:697:VAL:HG11	2:S:714:MET:HE2	2.00	0.43
2:S:845:LEU:O	2:S:849:VAL:HG13	2.17	0.43
2:S:677:PHE:HD1	2:S:677:PHE:HA	1.70	0.43
1:R:67:MET:HE1	2:S:929:PHE:CZ	2.54	0.43
2:S:585:ASN:O	2:S:604:GLY:HA2	2.18	0.43
2:S:852:VAL:O	2:S:856:ILE:HG13	2.19	0.43
1:R:42:LYS:C	1:R:52:LEU:HD12	2.43	0.43
1:R:19:LEU:HD23	1:R:19:LEU:HA	1.86	0.42
1:R:27:HIS:ND1	1:R:149:ARG:HD2	2.34	0.42
1:R:120:LEU:H	1:R:120:LEU:HD23	1.83	0.42
2:S:1033:TYR:HB2	2:S:1034:PRO:HD2	2.01	0.42
1:R:141:TYR:OH	1:R:143:GLU:HG2	2.20	0.42
2:S:601:ILE:HD11	2:S:612:ARG:CB	2.49	0.42
2:S:821:LEU:HD11	2:S:934:LEU:HD11	2.02	0.42
2:S:684:PRO:HA	2:S:687:LEU:HB2	2.02	0.42
1:R:37:GLU:O	1:R:41:ARG:HG3	2.20	0.42
2:S:690:LEU:HD21	2:S:721:VAL:HG11	2.00	0.42
1:R:73:ARG:NH2	2:S:879:ASN:O	2.53	0.41
2:S:754:PHE:CD1	2:S:980:CYS:HB2	2.56	0.41
2:S:831:LEU:HD12	2:S:831:LEU:HA	1.82	0.41
2:S:843:GLU:O	2:S:848:ARG:NH2	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:6:VAL:O	3:P:7:ALA:HB2	2.20	0.41
1:R:141:TYR:OH	1:R:143:GLU:OE1	2.33	0.41
2:S:828:THR:HG21	2:S:869:ASN:HD21	1.86	0.41
1:R:144:THR:HG22	1:R:151:GLY:C	2.46	0.41
2:S:572:ALA:C	2:S:574:VAL:H	2.29	0.41
2:S:576:ARG:HD3	2:S:651:PRO:HD3	2.03	0.41
1:R:71:TYR:O	1:R:74:THR:HB	2.21	0.40
2:S:642:SER:CA	2:S:713:ARG:HH22	2.34	0.40
2:S:627:PHE:O	2:S:631:TYR:HB3	2.21	0.40
2:S:824:MET:HE3	2:S:824:MET:HB3	1.92	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	R	164/186 (88%)	153 (93%)	11 (7%)	0	100	100
2	S	432/507 (85%)	415 (96%)	17 (4%)	0	100	100
3	P	6/8 (75%)	1 (17%)	1 (17%)	4 (67%)	0	0
All	All	602/701 (86%)	569 (94%)	29 (5%)	4 (1%)	18	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	P	5	SER
3	P	6	VAL
3	P	7	ALA
3	P	2	HIS

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	R	144/161 (89%)	137 (95%)	7 (5%)	22	54
2	S	414/469 (88%)	404 (98%)	10 (2%)	43	75
3	P	5/5 (100%)	4 (80%)	1 (20%)	1	4
All	All	563/635 (89%)	545 (97%)	18 (3%)	34	68

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	R	24	ILE
1	R	26	ASN
1	R	39	SER
1	R	41	ARG
1	R	68	ARG
1	R	88	LYS
1	R	128	ARG
2	S	615	TYR
2	S	675	LYS
2	S	677	PHE
2	S	678	ARG
2	S	682	ILE
2	S	726	MET
2	S	730	VAL
2	S	885	ARG
2	S	930	PHE
2	S	964	VAL
3	P	2	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	R	25	GLN
1	R	61	GLN
1	R	70	GLN

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Mol	Chain	Res	Type
2	S	700	HIS
2	S	800	GLN
2	S	936	ASN
2	S	972	GLN
3	P	2	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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
3	DAR	P	1	3	9,10,11	3.06	3 (33%)	5,11,13	3.16	1 (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
3	DAR	P	1	3	-	2/8/9/11	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	1	DAR	CZ-NE	7.50	1.47	1.33
3	P	1	DAR	CZ-NH2	3.67	1.45	1.32
3	P	1	DAR	CZ-NH1	-2.24	1.26	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	1	DAR	CG-CD-NE	7.06	132.00	112.20

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	P	1	DAR	NE-CD-CG-CB
3	P	1	DAR	CG-CD-NE-CZ

There are no ring outliers.

No monomer is involved in short contacts.

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	R	166/186 (89%)	0.02	2 (1%) 76 69	53, 73, 114, 146	1 (0%)
2	S	440/507 (86%)	0.12	6 (1%) 73 65	43, 81, 115, 136	19 (4%)
3	P	6/8 (75%)	0.91	0 100 100	127, 136, 182, 187	0
All	All	612/701 (87%)	0.10	8 (1%) 75 67	43, 79, 117, 187	20 (3%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	R	30	ASP	2.9
2	S	577	PHE	2.7
2	S	569	LEU	2.6
2	S	578	ALA	2.6
2	S	575	TYR	2.3
2	S	677	PHE	2.2
2	S	680	GLU	2.1
1	R	48	GLY	2.1

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
3	DAR	P	1	11/12	0.31	0.14	112,139,167,183	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.