



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 6, 2017 – 11:37 AM EST

PDB ID : 4H3K
Title : Crystal structure of a ternary complex of human symplekin NTD, human Ssu72 and a RNA polymerase II CTD peptide phosphorylated at Ser-2, Ser-5 and Ser-7
Authors : Xiang, K.; Tong, L.
Deposited on : unknown
Resolution : 2.00 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030345

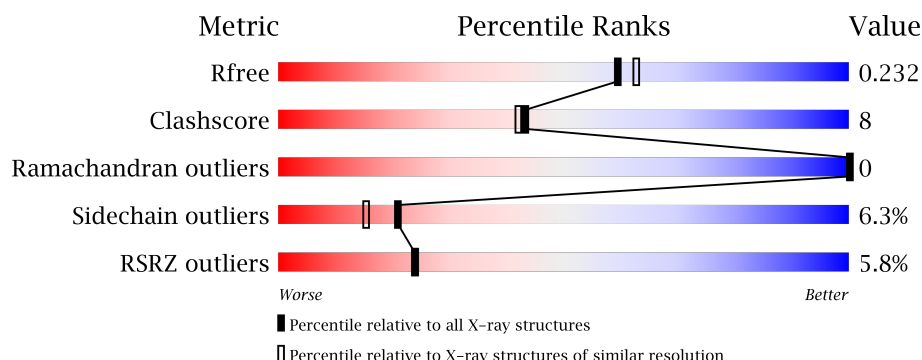
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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	351	<div> <div>5%</div> <div> <div></div> <div>79%</div> <div>9%</div> <div>10%</div> </div> </div>
1	D	351	<div> <div>8%</div> <div> <div></div> <div>75%</div> <div>11%</div> <div>11%</div> </div> </div>
2	B	214	<div> <div>4%</div> <div> <div></div> <div>68%</div> <div>17%</div> <div>12%</div> </div> </div>
2	E	214	<div> <div>3%</div> <div> <div></div> <div>70%</div> <div>16%</div> <div>11%</div> </div> </div>
3	F	10	<div> <div>10%</div> <div> <div></div> <div>30%</div> <div>40%</div> <div>30%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8247 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 Symplekin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	0	0
			2506	1596	433	464	13			
1	D	311	Total	C	N	O	S	0	0	0
			2461	1569	422	457	13			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	MET	-	EXPRESSION TAG	UNP Q92797
A	11	GLY	-	EXPRESSION TAG	UNP Q92797
A	12	SER	-	EXPRESSION TAG	UNP Q92797
A	13	SER	-	EXPRESSION TAG	UNP Q92797
A	14	HIS	-	EXPRESSION TAG	UNP Q92797
A	15	HIS	-	EXPRESSION TAG	UNP Q92797
A	16	HIS	-	EXPRESSION TAG	UNP Q92797
A	17	HIS	-	EXPRESSION TAG	UNP Q92797
A	18	HIS	-	EXPRESSION TAG	UNP Q92797
A	19	HIS	-	EXPRESSION TAG	UNP Q92797
A	20	SER	-	EXPRESSION TAG	UNP Q92797
A	21	SER	-	EXPRESSION TAG	UNP Q92797
A	22	GLY	-	EXPRESSION TAG	UNP Q92797
A	23	LEU	-	EXPRESSION TAG	UNP Q92797
A	24	VAL	-	EXPRESSION TAG	UNP Q92797
A	25	PRO	-	EXPRESSION TAG	UNP Q92797
A	26	ARG	-	EXPRESSION TAG	UNP Q92797
A	27	GLY	-	EXPRESSION TAG	UNP Q92797
A	28	SER	-	EXPRESSION TAG	UNP Q92797
A	29	HIS	-	EXPRESSION TAG	UNP Q92797
D	10	MET	-	EXPRESSION TAG	UNP Q92797
D	11	GLY	-	EXPRESSION TAG	UNP Q92797
D	12	SER	-	EXPRESSION TAG	UNP Q92797
D	13	SER	-	EXPRESSION TAG	UNP Q92797
D	14	HIS	-	EXPRESSION TAG	UNP Q92797

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Chain	Residue	Modelled	Actual	Comment	Reference
D	15	HIS	-	EXPRESSION TAG	UNP Q92797
D	16	HIS	-	EXPRESSION TAG	UNP Q92797
D	17	HIS	-	EXPRESSION TAG	UNP Q92797
D	18	HIS	-	EXPRESSION TAG	UNP Q92797
D	19	HIS	-	EXPRESSION TAG	UNP Q92797
D	20	SER	-	EXPRESSION TAG	UNP Q92797
D	21	SER	-	EXPRESSION TAG	UNP Q92797
D	22	GLY	-	EXPRESSION TAG	UNP Q92797
D	23	LEU	-	EXPRESSION TAG	UNP Q92797
D	24	VAL	-	EXPRESSION TAG	UNP Q92797
D	25	PRO	-	EXPRESSION TAG	UNP Q92797
D	26	ARG	-	EXPRESSION TAG	UNP Q92797
D	27	GLY	-	EXPRESSION TAG	UNP Q92797
D	28	SER	-	EXPRESSION TAG	UNP Q92797
D	29	HIS	-	EXPRESSION TAG	UNP Q92797

- Molecule 2 is a protein called RNA polymerase II subunit A C-terminal domain phosphatase SSU72.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	189	Total	C	N	O	S	0	0	0
			1546	960	272	303	11			
2	E	190	Total	C	N	O	S	0	0	0
			1553	965	273	304	11			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	EXPRESSION TAG	UNP Q9NP77
B	-18	GLY	-	EXPRESSION TAG	UNP Q9NP77
B	-17	SER	-	EXPRESSION TAG	UNP Q9NP77
B	-16	SER	-	EXPRESSION TAG	UNP Q9NP77
B	-15	HIS	-	EXPRESSION TAG	UNP Q9NP77
B	-14	HIS	-	EXPRESSION TAG	UNP Q9NP77
B	-13	HIS	-	EXPRESSION TAG	UNP Q9NP77
B	-12	HIS	-	EXPRESSION TAG	UNP Q9NP77
B	-11	HIS	-	EXPRESSION TAG	UNP Q9NP77
B	-10	HIS	-	EXPRESSION TAG	UNP Q9NP77
B	-9	SER	-	EXPRESSION TAG	UNP Q9NP77
B	-8	SER	-	EXPRESSION TAG	UNP Q9NP77
B	-7	GLY	-	EXPRESSION TAG	UNP Q9NP77
B	-6	LEU	-	EXPRESSION TAG	UNP Q9NP77

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	VAL	-	EXPRESSION TAG	UNP Q9NP77
B	-4	PRO	-	EXPRESSION TAG	UNP Q9NP77
B	-3	ARG	-	EXPRESSION TAG	UNP Q9NP77
B	-2	GLY	-	EXPRESSION TAG	UNP Q9NP77
B	-1	SER	-	EXPRESSION TAG	UNP Q9NP77
B	0	HIS	-	EXPRESSION TAG	UNP Q9NP77
B	12	SER	CYS	ENGINEERED MUTATION	UNP Q9NP77
E	-19	MET	-	EXPRESSION TAG	UNP Q9NP77
E	-18	GLY	-	EXPRESSION TAG	UNP Q9NP77
E	-17	SER	-	EXPRESSION TAG	UNP Q9NP77
E	-16	SER	-	EXPRESSION TAG	UNP Q9NP77
E	-15	HIS	-	EXPRESSION TAG	UNP Q9NP77
E	-14	HIS	-	EXPRESSION TAG	UNP Q9NP77
E	-13	HIS	-	EXPRESSION TAG	UNP Q9NP77
E	-12	HIS	-	EXPRESSION TAG	UNP Q9NP77
E	-11	HIS	-	EXPRESSION TAG	UNP Q9NP77
E	-10	HIS	-	EXPRESSION TAG	UNP Q9NP77
E	-9	SER	-	EXPRESSION TAG	UNP Q9NP77
E	-8	SER	-	EXPRESSION TAG	UNP Q9NP77
E	-7	GLY	-	EXPRESSION TAG	UNP Q9NP77
E	-6	LEU	-	EXPRESSION TAG	UNP Q9NP77
E	-5	VAL	-	EXPRESSION TAG	UNP Q9NP77
E	-4	PRO	-	EXPRESSION TAG	UNP Q9NP77
E	-3	ARG	-	EXPRESSION TAG	UNP Q9NP77
E	-2	GLY	-	EXPRESSION TAG	UNP Q9NP77
E	-1	SER	-	EXPRESSION TAG	UNP Q9NP77
E	0	HIS	-	EXPRESSION TAG	UNP Q9NP77
E	12	SER	CYS	ENGINEERED MUTATION	UNP Q9NP77

- Molecule 3 is a protein called Hexapeptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	7	Total	C	N	O	P	0	0	0
			63	32	7	21	3			

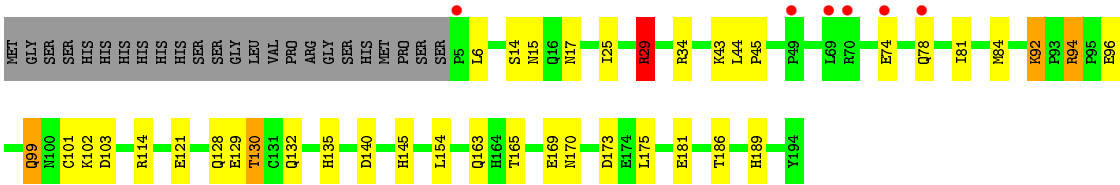
- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



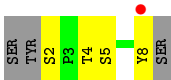
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	29	Total	O	0	0
			29	29		
5	B	29	Total	O	0	0
			29	29		
5	D	29	Total	O	0	0
			29	29		
5	E	26	Total	O	0	0
			26	26		



● Molecule 3: Hexapeptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.58 Å 97.36 Å 105.28 Å 90.00° 97.86° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 44.11 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.1 (50.00-2.00) 96.1 (44.11-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 2.00 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.197 , 0.233 0.198 , 0.232	Depositor DCC
R_{free} test set	4376 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 44.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8247	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, SEP

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	0.68	0/2544	0.71	2/3449 (0.1%)
1	D	0.69	0/2497	0.70	2/3385 (0.1%)
2	B	0.73	1/1574 (0.1%)	0.70	1/2124 (0.0%)
2	E	0.74	0/1582	0.81	3/2135 (0.1%)
3	F	0.52	0/33	0.42	0/41
All	All	0.70	1/8230 (0.0%)	0.73	8/11134 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	161	CYS	CB-SG	-9.59	1.66	1.82

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	29	ARG	NE-CZ-NH2	-9.86	115.37	120.30
1	A	315	LEU	CA-CB-CG	8.15	134.04	115.30
2	E	29	ARG	NE-CZ-NH1	7.41	124.00	120.30
1	D	218	ARG	NE-CZ-NH2	-7.05	116.78	120.30
1	D	160	MET	CG-SD-CE	-6.70	89.48	100.20
2	B	139	VAL	CB-CA-C	-6.33	99.37	111.40
1	A	245	HIS	CB-CA-C	5.58	121.57	110.40
2	E	140	ASP	CB-CG-OD1	5.46	123.22	118.30

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	2506	0	2610	29	1
1	D	2461	0	2566	29	1
2	B	1546	0	1493	39	0
2	E	1553	0	1501	38	0
3	F	63	0	42	4	0
4	B	5	0	0	0	0
5	A	29	0	0	0	0
5	B	29	0	0	4	0
5	D	29	0	0	0	0
5	E	26	0	0	2	0
All	All	8247	0	8212	125	1

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 (125) 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:E:165:THR:HG21	2:E:175:LEU:HD11	1.09	1.07
2:E:165:THR:HG21	2:E:175:LEU:CD1	1.93	0.98
2:E:99:GLN:HE21	2:E:99:GLN:H	1.09	0.98
2:B:135:HIS:HD2	2:B:189:HIS:CE1	1.84	0.95
2:B:99:GLN:HE21	2:B:99:GLN:H	1.13	0.94
2:E:135:HIS:HD2	2:E:189:HIS:CE1	1.86	0.93
2:B:60:THR:HG22	2:B:63:GLN:H	1.30	0.93
2:B:102:LYS:HD2	2:B:102:LYS:N	1.84	0.92
2:B:102:LYS:H	2:B:102:LYS:HD2	1.33	0.91
2:B:135:HIS:HD2	2:B:189:HIS:HE1	1.17	0.91
2:B:165:THR:HG21	2:B:175:LEU:HD11	1.53	0.91
2:E:165:THR:CG2	2:E:175:LEU:HD11	2.02	0.88
2:E:135:HIS:HD2	2:E:189:HIS:HE1	1.21	0.84
2:B:135:HIS:CD2	2:B:189:HIS:HE1	1.98	0.82
2:E:135:HIS:CD2	2:E:189:HIS:HE1	1.99	0.80
2:B:102:LYS:CD	2:B:102:LYS:H	1.95	0.78
2:B:15:ASN:HD22	2:B:96:GLU:H	1.31	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:15:ASN:HD21	2:E:92:LYS:NZ	1.82	0.77
1:A:144:MET:HB3	1:A:223:ILE:HD11	1.68	0.75
2:E:15:ASN:HD22	2:E:96:GLU:H	1.33	0.74
2:E:173:ASP:OD1	2:E:189:HIS:HD2	1.71	0.73
1:D:206:ARG:NE	2:E:129:GLU:HG2	2.04	0.71
2:E:99:GLN:HE21	2:E:99:GLN:N	1.86	0.71
2:B:130:THR:CG2	2:B:132:GLN:H	2.04	0.71
2:E:25:ILE:O	2:E:29:ARG:HG2	1.91	0.71
2:B:165:THR:CG2	2:B:175:LEU:HD11	2.21	0.70
1:D:177:ASP:H	2:E:170:ASN:HD22	1.39	0.70
2:B:102:LYS:CD	2:B:102:LYS:N	2.54	0.69
2:B:15:ASN:HD21	2:B:92:LYS:NZ	1.90	0.69
1:A:220:HIS:HD2	1:A:222:TYR:H	1.42	0.67
1:D:220:HIS:HD2	1:D:223:ILE:H	1.42	0.66
1:A:247:ALA:HB2	1:D:326:VAL:HG11	1.76	0.66
2:B:173:ASP:OD1	2:B:189:HIS:HD2	1.78	0.66
2:E:130:THR:HG23	2:E:132:GLN:H	1.61	0.65
1:D:220:HIS:CD2	1:D:223:ILE:H	2.15	0.65
1:A:312:PRO:O	1:A:315:LEU:HD13	1.98	0.64
2:E:135:HIS:HE1	2:E:169:GLU:OE2	1.81	0.64
1:A:220:HIS:CD2	1:A:222:TYR:H	2.16	0.64
2:E:45:PRO:HG2	2:E:81:ILE:HD11	1.78	0.63
1:A:339:MET:CB	1:A:340:PRO:HD3	2.29	0.63
1:A:177:ASP:H	2:B:170:ASN:HD22	1.46	0.62
1:A:246:PRO:HB3	1:D:332:GLN:HG3	1.82	0.62
1:D:150:ILE:HD12	1:D:222:TYR:CD2	2.35	0.61
2:B:66:ASN:O	2:B:70:ARG:HG3	2.00	0.61
2:B:15:ASN:HD21	2:B:92:LYS:HZ1	1.49	0.60
2:E:130:THR:CG2	2:E:132:GLN:H	2.14	0.60
2:B:130:THR:HG23	2:B:132:GLN:H	1.67	0.59
1:D:58:VAL:HG11	1:D:77:ILE:HD11	1.84	0.59
1:A:177:ASP:H	2:B:170:ASN:ND2	2.00	0.59
1:D:254:THR:HG21	5:E:217:HOH:O	2.03	0.59
1:D:318:GLN:O	1:D:322:THR:HB	2.03	0.59
1:A:339:MET:HB2	1:A:340:PRO:HD3	1.83	0.59
2:B:99:GLN:NE2	2:B:99:GLN:H	1.94	0.59
2:E:84:MET:HE3	2:E:145:HIS:NE2	2.17	0.58
1:D:220:HIS:HD2	1:D:222:TYR:H	1.51	0.57
1:A:285:LEU:HD23	1:A:285:LEU:H	1.70	0.57
1:A:59:GLN:HG3	1:A:63:ILE:HD12	1.87	0.57
2:B:130:THR:HG22	2:B:132:GLN:H	1.68	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:PRO:C	1:D:322:THR:HG21	2.25	0.56
1:A:254:THR:HG21	5:B:308:HOH:O	2.04	0.56
2:E:84:MET:CE	2:E:145:HIS:NE2	2.68	0.55
1:A:254:THR:CG2	5:B:308:HOH:O	2.53	0.55
2:B:165:THR:HB	2:B:171:GLU:OE1	2.05	0.55
2:E:135:HIS:CD2	2:E:189:HIS:CE1	2.75	0.55
1:D:220:HIS:CD2	1:D:222:TYR:H	2.24	0.55
1:A:198:MET:SD	1:A:269:MET:HG3	2.47	0.55
1:D:254:THR:CG2	5:E:217:HOH:O	2.54	0.54
2:E:34:ARG:NH1	2:E:103:ASP:OD2	2.41	0.54
2:E:74:GLU:O	2:E:78:GLN:HB2	2.07	0.54
2:B:13:SER:OG	2:B:114:ARG:NH2	2.40	0.54
1:D:177:ASP:H	2:E:170:ASN:ND2	2.04	0.54
1:D:239:LEU:O	1:D:243:MET:HG3	2.08	0.54
1:A:316:GLU:O	1:A:316:GLU:HG3	2.08	0.54
2:E:15:ASN:HD21	2:E:92:LYS:HZ1	1.56	0.53
2:E:15:ASN:HD21	2:E:92:LYS:HZ3	1.53	0.53
1:D:206:ARG:CZ	2:E:129:GLU:HG2	2.39	0.53
2:B:109:LEU:HD13	2:B:155:ILE:HG23	1.91	0.52
2:E:43:LYS:HB3	3:F:4:THR:HG22	1.93	0.51
2:B:60:THR:HG22	2:B:63:GLN:N	2.12	0.50
2:B:39:GLY:O	2:B:97:ARG:NH1	2.44	0.50
2:E:15:ASN:ND2	2:E:96:GLU:H	2.05	0.50
1:A:198:MET:SD	1:A:269:MET:CG	3.00	0.50
2:B:135:HIS:HE1	2:B:169:GLU:OE2	1.95	0.50
2:B:135:HIS:CD2	2:B:189:HIS:CE1	2.75	0.50
2:E:114:ARG:NH2	3:F:2:SEP:O1P	2.31	0.49
2:B:130:THR:CG2	2:B:132:GLN:HB3	2.42	0.49
1:A:220:HIS:CD2	1:A:223:ILE:H	2.31	0.49
2:B:31:PHE:HE1	2:B:163:GLN:HG3	1.79	0.48
2:E:45:PRO:HB3	3:F:8:TYR:HA	1.96	0.48
1:A:74:LEU:O	1:A:78:ILE:HG12	2.14	0.48
2:E:17:ASN:ND2	2:E:84:MET:HE2	2.29	0.47
1:D:282:HIS:HE1	1:D:327:ASP:OD2	1.98	0.47
2:B:34:ARG:NH1	2:B:103:ASP:OD2	2.48	0.47
1:D:185:LYS:NZ	2:E:128:GLN:HE22	2.13	0.47
2:E:6:LEU:H	2:E:163:GLN:HE21	1.63	0.47
2:E:14:SER:HB2	3:F:5:SEP:HB2	1.97	0.47
1:D:243:MET:SD	1:D:260:LEU:HD11	2.56	0.46
1:D:30:MET:HE2	1:D:35:ARG:HG2	1.99	0.45
2:E:101:CYS:C	2:E:102:LYS:HD2	2.37	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:92:LYS:HE3	2:E:94:ARG:O	2.17	0.45
2:B:6:LEU:N	5:B:323:HOH:O	2.51	0.44
1:D:118:LEU:HB2	1:D:164:MET:HE3	1.98	0.44
2:B:116:TYR:CD1	2:B:138:ASN:HB2	2.53	0.44
1:A:25:PRO:O	1:A:26:ARG:HG2	2.18	0.44
1:A:58:VAL:HG11	1:A:77:ILE:HD11	2.00	0.44
1:D:250:SER:O	1:D:254:THR:HG23	2.18	0.43
2:B:165:THR:HG21	2:B:175:LEU:CD1	2.38	0.43
1:A:250:SER:O	1:A:254:THR:HG23	2.19	0.43
1:A:118:LEU:HB2	1:A:164:MET:HE3	2.01	0.42
1:A:220:HIS:HD2	1:A:223:ILE:H	1.65	0.42
1:D:113:ASN:O	1:D:116:MET:HG3	2.19	0.42
1:D:150:ILE:HG23	1:D:154:GLN:HB3	2.02	0.42
1:A:316:GLU:HG2	1:A:317:PHE:CZ	2.55	0.42
2:B:17:ASN:ND2	5:B:321:HOH:O	2.44	0.42
1:D:59:GLN:HG3	1:D:63:ILE:HD12	2.02	0.42
2:B:60:THR:CG2	2:B:62:ASP:HB3	2.50	0.42
1:A:31:THR:OG1	1:A:34:GLU:HG3	2.19	0.41
2:B:15:ASN:ND2	2:B:96:GLU:H	2.08	0.41
1:D:114:LEU:HD13	1:D:133:MET:SD	2.61	0.41
1:D:109:LYS:HG2	1:D:109:LYS:H	1.62	0.41
1:A:118:LEU:HD22	1:A:164:MET:HE2	2.03	0.41
1:A:144:MET:HB3	1:A:223:ILE:CD1	2.46	0.40
2:B:60:THR:HG23	2:B:62:ASP:H	1.87	0.40
1:D:41:ASN:N	1:D:41:ASN:HD22	2.20	0.40
2:E:25:ILE:O	2:E:29:ARG:CG	2.66	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:GLU:OE2	1:D:218:ARG:NH2[2_455]	2.18	0.02

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	315/351 (90%)	310 (98%)	5 (2%)	0	100	100
1	D	309/351 (88%)	306 (99%)	3 (1%)	0	100	100
2	B	187/214 (87%)	184 (98%)	3 (2%)	0	100	100
2	E	188/214 (88%)	184 (98%)	4 (2%)	0	100	100
3	F	3/10 (30%)	3 (100%)	0	0	100	100
All	All	1002/1140 (88%)	987 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	283/315 (90%)	268 (95%)	15 (5%)	26	21
1	D	278/315 (88%)	258 (93%)	20 (7%)	17	11
2	B	177/199 (89%)	164 (93%)	13 (7%)	16	11
2	E	178/199 (89%)	168 (94%)	10 (6%)	25	19
3	F	4/7 (57%)	4 (100%)	0	100	100
All	All	920/1035 (89%)	862 (94%)	58 (6%)	21	15

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ARG
1	A	71	ASP
1	A	114	LEU
1	A	116	MET
1	A	148	ARG
1	A	149	VAL
1	A	198	MET

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Mol	Chain	Res	Type
1	A	209	HIS
1	A	213	LEU
1	A	215	ARG
1	A	245	HIS
1	A	254	THR
1	A	269	MET
1	A	285	LEU
1	A	312	PRO
2	B	29	ARG
2	B	43	LYS
2	B	44	LEU
2	B	60	THR
2	B	68	LEU
2	B	73	LYS
2	B	78	GLN
2	B	99	GLN
2	B	102	LYS
2	B	103	ASP
2	B	130	THR
2	B	163	GLN
2	B	186	THR
1	D	35	ARG
1	D	41	ASN
1	D	65	LYS
1	D	72	ASN
1	D	81	GLN
1	D	102	ARG
1	D	109	LYS
1	D	116	MET
1	D	148	ARG
1	D	160	MET
1	D	198	MET
1	D	205	ARG
1	D	211	ILE
1	D	213	LEU
1	D	215	ARG
1	D	254	THR
1	D	285	LEU
1	D	291	LYS
1	D	307	SER
1	D	322	THR
2	E	29	ARG

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Mol	Chain	Res	Type
2	E	44	LEU
2	E	92	LYS
2	E	94	ARG
2	E	99	GLN
2	E	121	GLU
2	E	130	THR
2	E	154	LEU
2	E	181	GLU
2	E	186	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	59	GLN
1	A	81	GLN
1	A	113	ASN
1	A	142	GLN
1	A	207	GLN
1	A	220	HIS
1	A	282	HIS
2	B	15	ASN
2	B	17	ASN
2	B	99	GLN
2	B	135	HIS
2	B	164	HIS
2	B	170	ASN
2	B	189	HIS
1	D	41	ASN
1	D	59	GLN
1	D	81	GLN
1	D	113	ASN
1	D	220	HIS
1	D	282	HIS
1	D	332	GLN
2	E	15	ASN
2	E	17	ASN
2	E	99	GLN
2	E	100	ASN
2	E	128	GLN
2	E	135	HIS
2	E	163	GLN

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Mol	Chain	Res	Type
2	E	170	ASN
2	E	189	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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
3	SEP	F	2	3	9,9,10	1.96	2 (22%)	9,12,14	1.76	3 (33%)
3	SEP	F	5	3	9,9,10	1.79	2 (22%)	9,12,14	1.02	0
3	SEP	F	7	3	9,9,10	1.54	2 (22%)	9,12,14	1.41	1 (11%)

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	SEP	F	2	3	-	0/5/8/10	0/0/0/0
3	SEP	F	5	3	-	0/5/8/10	0/0/0/0
3	SEP	F	7	3	-	0/5/8/10	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	7	SEP	P-O3P	2.07	1.63	1.54
3	F	5	SEP	P-O1P	3.11	1.61	1.50
3	F	7	SEP	P-O1P	3.37	1.62	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	2	SEP	P-O1P	3.39	1.62	1.50
3	F	5	SEP	CA-C	3.39	1.54	1.50
3	F	2	SEP	CA-C	3.56	1.54	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	7	SEP	P-OG-CB	-3.41	108.91	118.30
3	F	2	SEP	O-C-CA	-2.05	119.36	125.02
3	F	2	SEP	OG-P-O1P	2.46	113.36	106.47
3	F	2	SEP	OG-CB-CA	3.28	111.40	108.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	2	SEP	1	0
3	F	5	SEP	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand 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	$\# Z > 2$
4	PO4	B	201	-	4,4,4	0.62	0	6,6,6	0.26	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
4	PO4	B	201	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	317/351 (90%)	0.12	16 (5%)	30 30	17, 34, 52, 65	0
1	D	311/351 (88%)	0.25	28 (9%)	10 10	17, 32, 74, 94	0
2	B	189/214 (88%)	0.20	8 (4%)	37 37	18, 33, 56, 66	0
2	E	190/214 (88%)	0.15	6 (3%)	48 48	18, 30, 53, 64	0
3	F	4/10 (40%)	1.38	1 (25%)	1 1	43, 44, 52, 58	0
All	All	1011/1140 (88%)	0.19	59 (5%)	24 24	17, 33, 58, 94	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	45	LEU	7.2
1	D	30	MET	6.6
1	A	148	ARG	4.9
1	D	148	ARG	4.6
1	D	31	THR	4.6
3	F	8	TYR	4.5
1	D	65	LYS	4.5
1	A	24	VAL	4.5
1	A	45	LEU	4.2
1	D	36	VAL	4.1
1	D	38	ASP	4.0
1	D	64	ASN	3.7
1	D	47	THR	3.7
1	D	40	LEU	3.5
1	A	47	THR	3.5
1	D	57	GLN	3.5
1	D	54	VAL	3.4
2	E	74	GLU	3.4
1	D	69	LEU	3.3
1	D	61	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	285	LEU	3.2
2	B	74	GLU	3.2
1	D	41	ASN	3.2
2	E	49	PRO	3.2
1	D	205	ARG	3.1
1	D	35	ARG	3.1
2	B	70	ARG	2.9
1	A	285	LEU	2.9
1	D	68	THR	2.8
2	B	69	LEU	2.8
1	A	246	PRO	2.8
1	D	66	ASP	2.8
1	A	27	GLY	2.7
1	D	46	ILE	2.7
1	D	42	GLN	2.7
2	E	69	LEU	2.6
1	A	25	PRO	2.6
1	D	53	THR	2.5
2	E	5	PRO	2.5
1	A	26	ARG	2.5
2	B	78	GLN	2.5
1	A	332	GLN	2.4
1	D	37	VAL	2.4
1	D	149	VAL	2.4
2	E	70	ARG	2.4
1	A	54	VAL	2.4
1	A	245	HIS	2.4
2	B	73	LYS	2.4
1	D	44	ALA	2.3
1	A	222	TYR	2.3
2	B	75	LEU	2.3
1	A	149	VAL	2.2
1	D	222	TYR	2.2
1	A	50	SER	2.1
1	D	60	GLU	2.1
1	A	205	ARG	2.1
2	B	83	HIS	2.1
2	B	166	GLU	2.0
2	E	78	GLN	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
3	SEP	F	7	10/11	0.72	0.20	-	48,52,64,66	0
3	SEP	F	5	10/11	0.99	0.10	-	25,33,39,40	0
3	SEP	F	2	10/11	0.74	0.30	-	55,57,69,70	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	PO4	B	201	5/5	0.99	0.11	0.06	30,30,33,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.