



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:11 AM UTC

PDB ID : 7YYL / pdb_00007yyL
Title : nucleotide-free DCCP:DCCP-R complex
Authors : Jeoung, J.-H.; Dobbek, H.
Deposited on : 2022-02-18
Resolution : 3.25 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

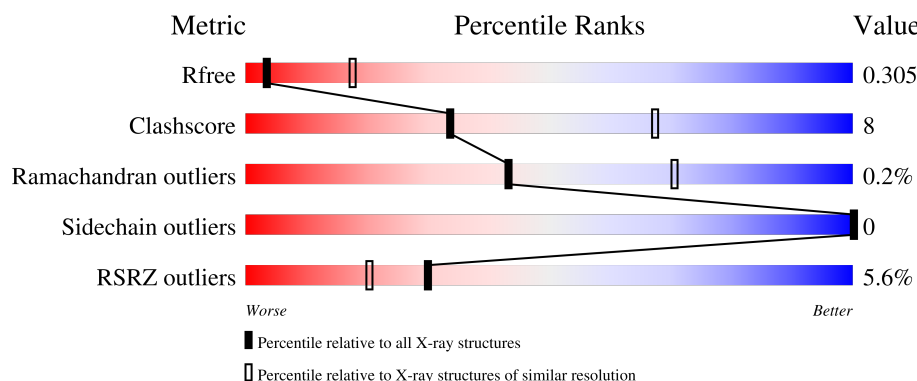
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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.



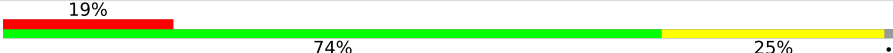
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1605 (3.30-3.22)
Clashscore	190562	1660 (3.30-3.22)
Ramachandran outliers	187476	1630 (3.30-3.22)
Sidechain outliers	187428	1629 (3.30-3.22)
RSRZ outliers	180081	1605 (3.30-3.22)

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	422	2% 80% 20%
1	B	422	2% 83% 16% .
1	C	422	% 82% 18%
1	D	422	3% 82% 17%
2	E	243	3% 85% 15%

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Mol	Chain	Length	Quality of chain
2	F	243	
2	G	243	
2	H	243	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 20929 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 Dehydratase family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	421	Total	C	N	O	S	1	1	0
			3371	2154	578	621	18			
1	D	420	Total	C	N	O	S	0	3	0
			3377	2158	579	623	17			
1	A	421	Total	C	N	O	S	0	1	0
			3371	2154	578	621	18			
1	B	419	Total	C	N	O	S	0	1	0
			3357	2146	578	616	17			

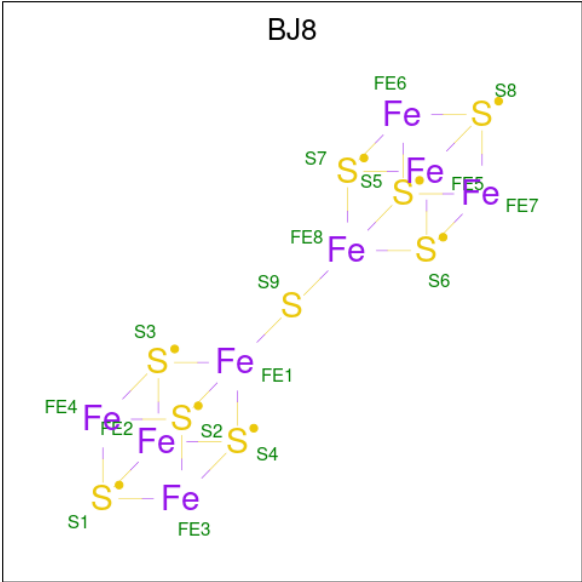
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	GLY	-	expression tag	UNP Q3AET9
D	0	GLY	-	expression tag	UNP Q3AET9
A	0	GLY	-	expression tag	UNP Q3AET9
B	0	GLY	-	expression tag	UNP Q3AET9

- Molecule 2 is a protein called Putative CoA-substrate-specific enzyme activase.

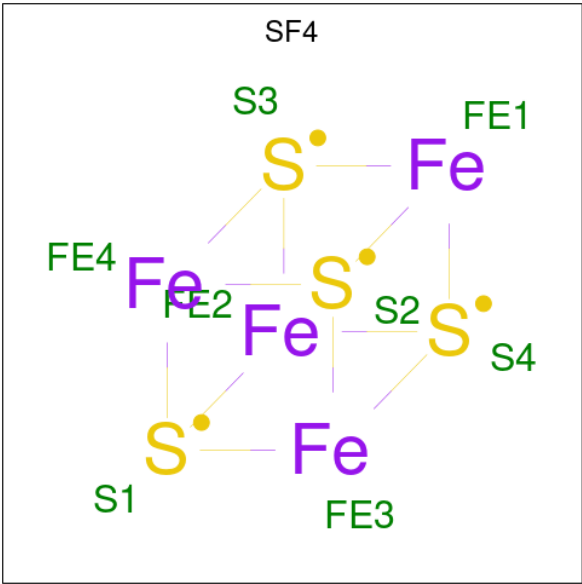
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	243	Total	C	N	O	S	0	0	0
			1848	1177	318	345	8			
2	F	241	Total	C	N	O	S	0	0	0
			1829	1163	316	343	7			
2	G	243	Total	C	N	O	S	0	0	0
			1848	1177	318	345	8			
2	H	241	Total	C	N	O	S	0	0	0
			1829	1163	316	343	7			

- Molecule 3 is Double cubane cluster (CCD ID: BJ8) (formula: Fe₈S₉) (labeled as "Ligand of Interest" by depositor).



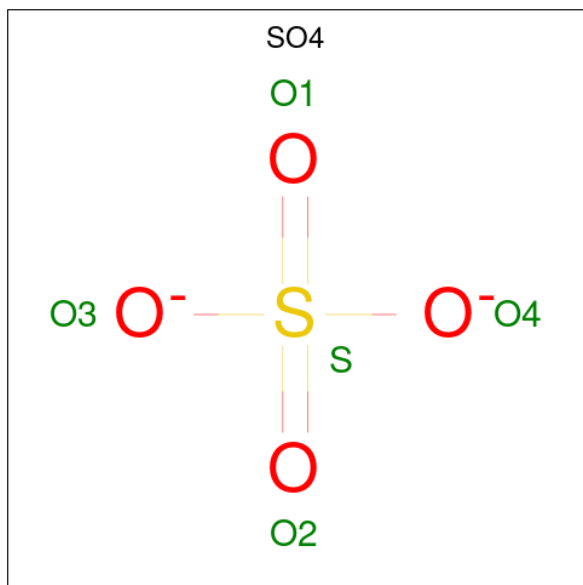
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	Fe	S	0	0
			17	8	9		
3	D	1	Total	Fe	S	0	0
			17	8	9		
3	A	1	Total	Fe	S	0	0
			17	8	9		
3	B	1	Total	Fe	S	0	0
			17	8	9		

- Molecule 4 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	E	1	Total	Fe	S	0	0
			8	4	4		
4	G	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

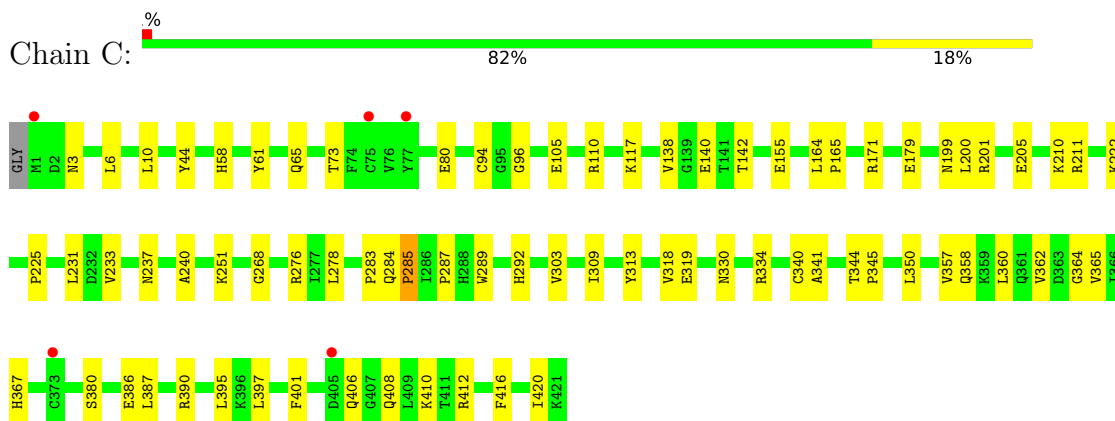


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	E	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	G	1	Total	O	S	0	0
			5	4	1		

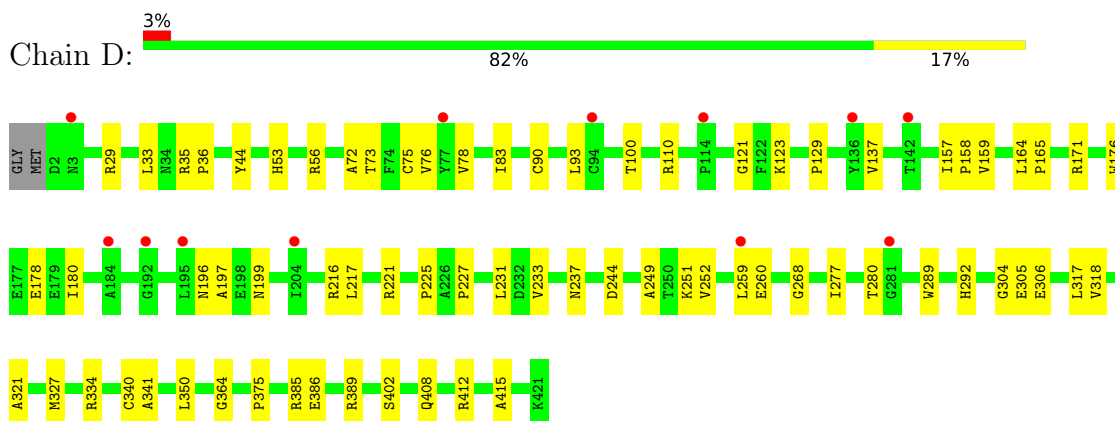
3 Residue-property plots [i](#)

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.

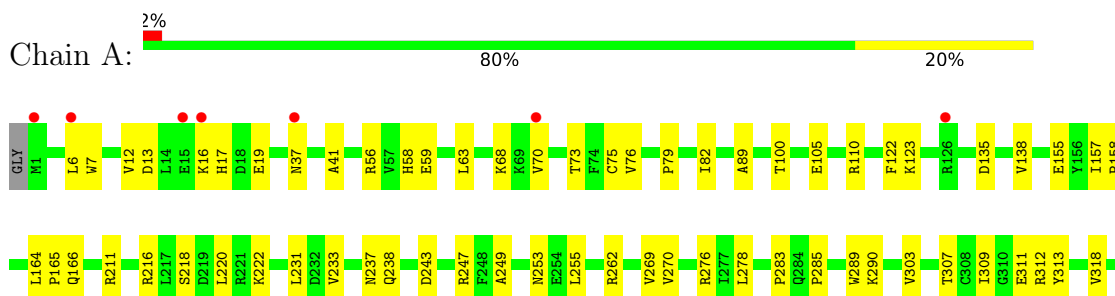
• Molecule 1: Dehydratase family protein



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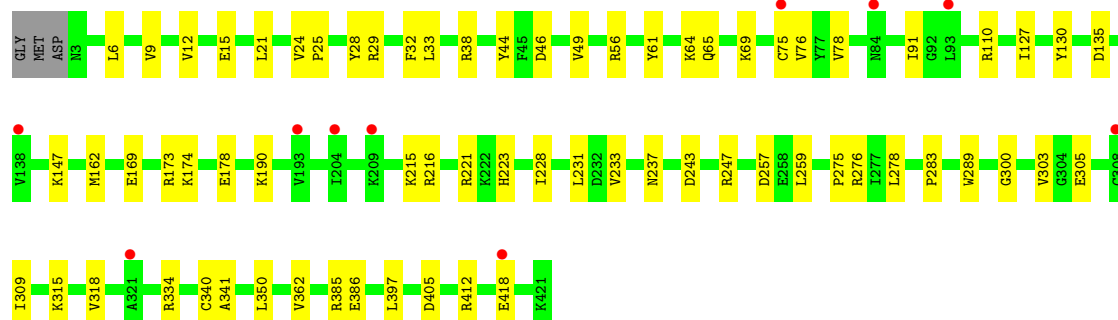
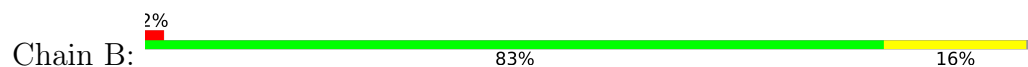


• Molecule 1: Dehydratase family protein

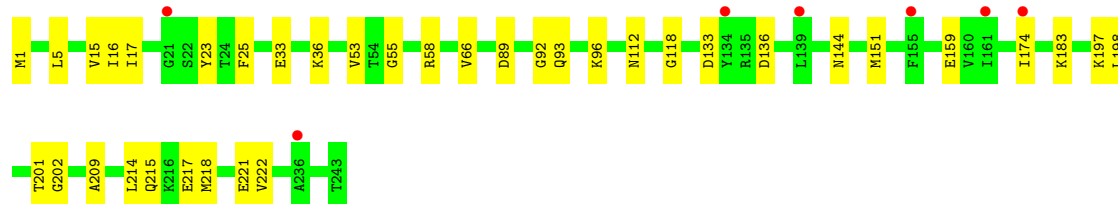
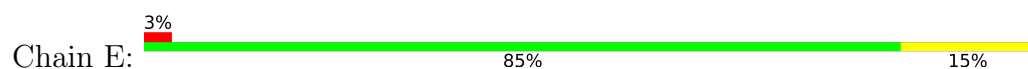




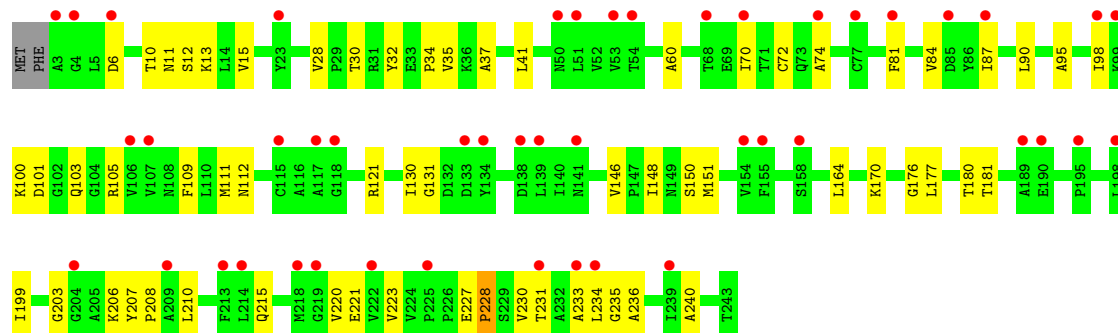
- Molecule 1: Dehydratase family protein



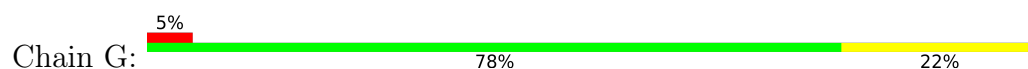
- Molecule 2: Putative CoA-substrate-specific enzyme activase



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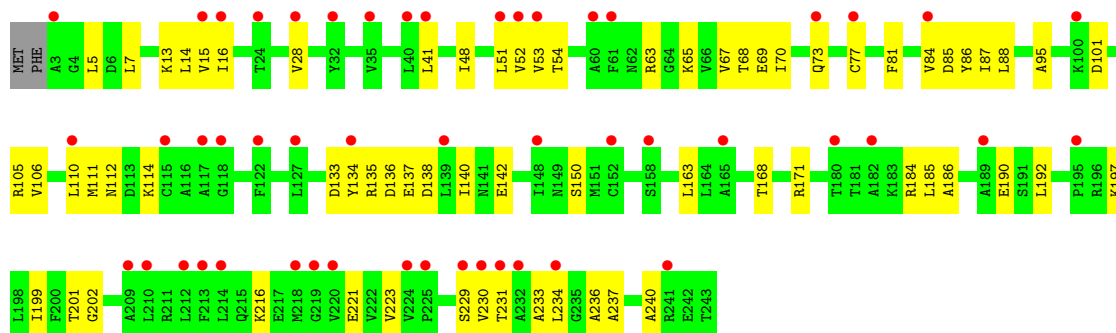


- Molecule 2: Putative CoA-substrate-specific enzyme activase





- Molecule 2: Putative CoA-substrate-specific enzyme activase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	81.67Å 83.33Å 124.38Å 99.72° 95.62° 94.02°	Depositor
Resolution (Å)	46.93 – 3.25 46.93 – 3.25	Depositor EDS
% Data completeness (in resolution range)	90.0 (46.93-3.25) 90.0 (46.93-3.25)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 3.25Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.246 , 0.303 0.248 , 0.305	Depositor DCC
R_{free} test set	2100 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	73.1	Xtriage
Anisotropy	0.141	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 80.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.000 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	20929	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.00% 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: BJ8, SF4, SO4

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	0.13	0/3443	0.30	0/4650
1	B	0.13	0/3429	0.33	0/4631
1	C	0.13	0/3443	0.33	0/4650
1	D	0.11	0/3455	0.30	0/4666
2	E	0.12	0/1874	0.32	0/2528
2	F	0.15	0/1854	0.36	0/2502
2	G	0.11	0/1874	0.32	0/2528
2	H	0.16	0/1854	0.42	0/2502
All	All	0.13	0/21226	0.33	0/28657

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 [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	3371	0	3401	53	0
1	B	3357	0	3390	48	0
1	C	3371	0	3401	49	0
1	D	3377	0	3406	46	0
2	E	1848	0	1919	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1829	0	1898	40	0
2	G	1848	0	1919	37	0
2	H	1829	0	1898	47	0
3	A	17	0	0	0	0
3	B	17	0	0	0	0
3	C	17	0	0	0	0
3	D	17	0	0	1	0
4	E	8	0	0	0	0
4	G	8	0	0	0	0
5	E	5	0	0	1	0
5	F	5	0	0	0	0
5	G	5	0	0	1	0
All	All	20929	0	21232	325	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (325) 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:H:69:GLU:O	2:H:73:GLN:OE1	1.71	1.06
1:A:406:GLN:O	1:A:410:LYS:HB2	1.74	0.87
1:C:406:GLN:O	1:C:410:LYS:HB2	1.74	0.87
1:A:233:VAL:O	1:A:237:ASN:ND2	2.08	0.86
2:F:148:ILE:HG13	2:F:177:LEU:HD22	1.65	0.78
1:B:412:ARG:NH1	2:H:112:ASN:O	2.17	0.77
2:E:214:LEU:O	2:E:218:MET:HB2	1.83	0.77
1:B:110:ARG:HH22	2:G:116:ALA:HB2	1.50	0.77
2:G:33:GLU:HB3	2:G:36:LYS:HD3	1.70	0.74
1:D:157:ILE:HG13	1:D:158:PRO:HD2	1.70	0.73
1:D:415:ALA:HB1	2:F:70:ILE:HD11	1.69	0.73
2:G:84:VAL:O	2:G:100:LYS:NZ	2.24	0.71
1:A:233:VAL:HG12	1:A:237:ASN:HD21	1.56	0.70
1:A:211:ARG:NH2	1:A:313:TYR:O	2.24	0.70
1:D:340:CYS:SG	1:D:341:ALA:N	2.65	0.70
1:D:412:ARG:NH1	2:F:112:ASN:O	2.25	0.70
1:B:340:CYS:SG	1:B:341:ALA:N	2.65	0.69
2:F:12:SER:HB2	2:F:28:VAL:HG13	1.74	0.69
1:A:231:LEU:HD11	1:A:289:TRP:HB3	1.74	0.69
1:B:350:LEU:HD21	1:B:386:GLU:HG2	1.74	0.69
2:E:58:ARG:HG3	2:E:66:VAL:HG23	1.72	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:LEU:HD21	1:A:386:GLU:HG2	1.75	0.69
2:H:13:LYS:HD3	2:H:230:VAL:HG22	1.74	0.69
1:D:231:LEU:HD11	1:D:289:TRP:HB3	1.76	0.68
2:F:13:LYS:HB3	2:F:230:VAL:HG12	1.76	0.68
2:H:15:VAL:HG21	2:H:234:LEU:HB3	1.75	0.68
1:C:350:LEU:HD21	1:C:386:GLU:HG2	1.76	0.67
1:D:164:LEU:HD12	1:D:165:PRO:HD2	1.75	0.67
1:B:69:LYS:NZ	1:B:190:LYS:O	2.25	0.66
2:G:73:GLN:NE2	2:G:89:ASP:OD2	2.23	0.66
2:F:28:VAL:HG11	2:F:41:LEU:HG	1.76	0.66
2:F:37:ALA:O	2:F:41:LEU:HB2	1.95	0.65
1:C:155:GLU:OE2	1:D:385:ARG:NH1	2.29	0.65
1:D:386:GLU:OE1	1:D:389:ARG:NH2	2.30	0.65
1:B:75:CYS:SG	1:B:76:VAL:N	2.69	0.65
2:F:164:LEU:HD22	2:F:170:LYS:HE3	1.79	0.65
2:H:14:LEU:HB2	2:H:41:LEU:HD13	1.77	0.64
1:B:283:PRO:HB3	1:B:309:ILE:HG13	1.79	0.64
2:G:35:VAL:HG23	2:G:36:LYS:HD2	1.79	0.64
1:A:307:THR:HA	1:A:312:ARG:HH21	1.62	0.64
2:E:214:LEU:O	2:E:218:MET:CB	2.47	0.63
2:H:70:ILE:HG22	2:H:106:VAL:HG11	1.79	0.63
1:D:75:CYS:SG	1:D:76:VAL:N	2.72	0.63
2:G:202:GLY:O	2:G:228:PRO:HB2	1.99	0.63
1:D:233:VAL:O	1:D:237:ASN:ND2	2.32	0.63
1:C:233:VAL:O	1:C:237:ASN:ND2	2.33	0.62
1:D:110:ARG:NH1	2:E:151:MET:O	2.33	0.61
1:C:164:LEU:HD12	1:C:165:PRO:HD2	1.81	0.61
2:E:33:GLU:OE1	2:E:36:LYS:NZ	2.28	0.61
2:F:103:GLN:OE1	2:F:105:ARG:NH2	2.33	0.61
1:A:70:VAL:HG23	1:A:89:ALA:HB3	1.81	0.60
2:H:77:CYS:HB2	2:H:87:ILE:HD12	1.83	0.60
1:A:41:ALA:HB3	1:A:255:LEU:HD13	1.82	0.60
1:B:174:LYS:NZ	1:B:178:GLU:OE2	2.33	0.60
1:A:75:CYS:SG	1:A:76:VAL:N	2.75	0.60
1:B:233:VAL:O	1:B:237:ASN:ND2	2.34	0.59
1:C:231:LEU:HD11	1:C:289:TRP:HB3	1.85	0.59
1:C:171:ARG:HH22	1:D:178:GLU:HB3	1.67	0.59
2:F:215:GLN:NE2	2:F:220:VAL:O	2.36	0.59
2:H:197:LYS:HE2	2:H:221:GLU:HB3	1.85	0.59
1:B:91:ILE:HD11	1:B:130:TYR:HB3	1.84	0.58
1:A:63:LEU:HD11	1:A:68:LYS:HD2	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:5:LEU:HD23	2:H:53:VAL:HG23	1.84	0.58
1:D:216:ARG:NH2	1:D:260:GLU:OE2	2.35	0.58
1:C:276:ARG:HE	1:C:303:VAL:HG11	1.69	0.58
1:A:105:GLU:OE1	1:A:110:ARG:NH2	2.36	0.58
1:C:80:GLU:OE2	1:C:210:LYS:NZ	2.36	0.58
1:C:110:ARG:NH1	2:F:151:MET:O	2.37	0.58
1:D:29:ARG:HA	1:D:33:LEU:HB2	1.86	0.58
2:E:89:ASP:OD2	2:E:96:LYS:NZ	2.36	0.58
1:C:386:GLU:OE2	1:C:390:ARG:NE	2.36	0.57
2:H:16:ILE:HD11	2:H:48:ILE:HD11	1.86	0.57
2:F:72:CYS:SG	2:F:236:ALA:N	2.78	0.57
1:A:222:LYS:HB3	1:A:360:LEU:HD23	1.87	0.57
1:C:179:GLU:OE2	1:D:171:ARG:NH2	2.38	0.57
2:H:52:VAL:HB	2:H:236:ALA:HB1	1.87	0.56
1:C:365:VAL:HG21	1:C:387:LEU:HD23	1.87	0.56
2:E:201:THR:OG1	2:E:202:GLY:N	2.37	0.56
2:F:11:ASN:HA	2:F:30:THR:H	1.71	0.56
1:D:72:ALA:HB1	1:D:93:LEU:HD21	1.86	0.56
2:F:228:PRO:HA	2:F:231:THR:OG1	2.06	0.56
2:H:63:ARG:O	2:H:63:ARG:HG2	2.05	0.56
2:F:72:CYS:O	2:F:235:GLY:HA3	2.06	0.55
1:A:73:THR:HG22	1:A:138:VAL:HB	1.87	0.55
1:C:201:ARG:O	1:C:205:GLU:HG2	2.07	0.55
1:A:16:LYS:HA	1:A:19:GLU:HB2	1.87	0.55
1:C:357:VAL:HG23	1:C:362:VAL:HB	1.87	0.55
1:C:222:LYS:HB3	1:C:360:LEU:HD23	1.89	0.55
2:G:87:ILE:HD11	2:G:100:LYS:HD3	1.89	0.55
2:H:137:GLU:HA	2:H:140:ILE:HD13	1.89	0.55
2:E:58:ARG:HG3	2:E:66:VAL:CG2	2.37	0.55
2:G:227:GLU:HB3	2:G:230:VAL:HG22	1.87	0.55
1:D:227:PRO:HB2	1:D:259:LEU:HD22	1.89	0.55
2:G:1:MET:N	2:G:17:ILE:O	2.40	0.54
1:C:225:PRO:HB2	1:C:268:GLY:HA3	1.89	0.54
1:A:249:ALA:O	1:A:253:ASN:ND2	2.39	0.54
2:H:88:LEU:HD21	2:H:185:LEU:HD11	1.89	0.54
1:B:6:LEU:O	1:B:9:VAL:HG12	2.07	0.54
2:E:92:GLY:N	5:E:302:SO4:O4	2.38	0.53
2:F:150:SER:HB3	2:F:180:THR:HG23	1.89	0.53
1:A:278:LEU:HD11	1:A:356:MET:HE2	1.90	0.53
1:D:318:VAL:HG22	1:D:334:ARG:HD3	1.91	0.53
2:F:35:VAL:HG22	2:F:60:ALA:HB1	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:VAL:HA	1:B:162:MET:HE1	1.90	0.53
1:B:405:ASP:OD2	2:H:114:LYS:NZ	2.31	0.53
2:F:181:THR:HB	2:F:210:LEU:HD11	1.91	0.52
1:D:277:ILE:HA	1:D:364:GLY:O	2.10	0.52
2:F:6:ASP:HB2	2:F:13:LYS:HB2	1.92	0.52
1:C:390:ARG:NH1	2:G:62:ASN:O	2.43	0.52
1:B:318:VAL:HG22	1:B:334:ARG:HD3	1.92	0.52
1:B:44:TYR:HE1	1:B:247:ARG:HE	1.56	0.52
1:B:223:HIS:O	1:B:276:ARG:NH1	2.42	0.52
1:D:233:VAL:HG12	1:D:237:ASN:HD21	1.74	0.51
1:D:292:HIS:NE2	1:D:305:GLU:OE2	2.37	0.51
1:B:29:ARG:HA	1:B:33:LEU:HD12	1.92	0.51
2:G:123:LEU:HD21	2:G:177:LEU:HD22	1.92	0.51
2:H:134:TYR:OH	2:H:171:ARG:HD2	2.11	0.51
1:B:215:LYS:NZ	1:B:315:LYS:O	2.31	0.51
1:C:412:ARG:NH2	2:E:112:ASN:O	2.45	0.50
2:F:98:ILE:HG22	2:F:109:PHE:HB2	1.93	0.50
2:G:211:ARG:NH2	2:G:222:VAL:O	2.45	0.50
1:C:6:LEU:HD22	1:C:58:HIS:CD2	2.46	0.50
2:G:13:LYS:NZ	5:G:302:SO4:O4	2.45	0.50
1:A:238:GLN:HE21	1:A:285:PRO:C	2.19	0.50
1:B:221:ARG:HH22	1:B:305:GLU:CD	2.19	0.50
2:G:56:TYR:OH	2:G:94:ASP:OD2	2.15	0.50
2:F:81:PHE:HZ	2:F:223:VAL:HG11	1.76	0.50
1:A:7:TRP:HB3	1:A:12:VAL:HG11	1.92	0.50
2:H:186:ALA:O	2:H:190:GLU:HG2	2.12	0.50
1:C:278:LEU:HB3	1:C:365:VAL:HG12	1.92	0.50
1:A:37:ASN:OD1	1:A:262:ARG:NH2	2.45	0.50
1:D:123:LYS:O	1:D:123:LYS:NZ	2.32	0.50
1:A:408:GLN:O	1:A:412:ARG:HD3	2.11	0.50
1:C:278:LEU:HB2	1:C:362:VAL:HG21	1.94	0.50
1:C:140:GLU:HG2	1:C:142:THR:HG23	1.93	0.49
1:C:416:PHE:O	1:C:420:ILE:HG12	2.13	0.49
1:D:44:TYR:CD2	1:D:251:LYS:HG2	2.46	0.49
2:F:236:ALA:O	2:F:240:ALA:N	2.44	0.49
1:B:12:VAL:HG13	1:B:127:ILE:HA	1.94	0.49
1:B:56:ARG:HE	1:B:243:ASP:HA	1.78	0.49
2:G:39:GLU:CD	2:G:63:ARG:HH12	2.20	0.49
1:C:3:ASN:OD1	1:C:58:HIS:NE2	2.44	0.49
2:E:93:GLN:HB2	2:E:118:GLY:H	1.77	0.49
2:F:221:GLU:HG2	2:F:223:VAL:HG22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:ASP:HB3	1:A:16:LYS:CE	2.42	0.49
1:B:278:LEU:HB2	1:B:362:VAL:HG21	1.95	0.49
2:G:138:ASP:O	2:G:171:ARG:NH1	2.43	0.49
1:D:197:ALA:HA	1:D:327:MET:HE1	1.94	0.49
2:H:231:THR:O	2:H:234:LEU:HG	2.13	0.49
1:A:395:LEU:HD22	1:A:416:PHE:HB2	1.94	0.49
2:H:95:ALA:N	2:H:112:ASN:OD1	2.44	0.49
1:D:225:PRO:HB2	1:D:268:GLY:HA3	1.94	0.49
2:F:101:ASP:OD1	2:F:101:ASP:N	2.46	0.48
2:F:130:ILE:HG13	2:F:131:GLY:H	1.78	0.48
1:B:110:ARG:NH2	2:G:113:ASP:OD2	2.43	0.48
1:A:13:ASP:HB3	1:A:16:LYS:HE3	1.95	0.48
2:H:70:ILE:HA	2:H:73:GLN:HE22	1.78	0.48
1:D:56:ARG:HH21	1:D:244:ASP:HA	1.79	0.48
1:A:312:ARG:NH1	1:A:376:TYR:OH	2.42	0.48
1:D:83:ILE:HG21	1:D:90:CYS:HB2	1.96	0.48
1:A:318:VAL:HG22	1:A:334:ARG:HD3	1.95	0.48
2:F:15:VAL:HG11	2:F:234:LEU:HD12	1.94	0.48
1:C:340:CYS:SG	1:C:341:ALA:N	2.87	0.48
1:C:284:GLN:O	1:C:309:ILE:HD12	2.14	0.48
1:A:70:VAL:HG12	1:A:135:ASP:H	1.79	0.48
1:A:389:ARG:O	1:A:389:ARG:NE	2.45	0.48
2:G:130:ILE:HD11	2:G:164:LEU:HD11	1.96	0.48
1:D:321:ALA:HB3	1:D:327:MET:HG3	1.96	0.48
2:G:17:ILE:HG12	2:G:23:TYR:HB3	1.96	0.48
2:G:146:VAL:O	2:G:176:GLY:HA3	2.14	0.47
2:E:55:GLY:O	2:E:58:ARG:NE	2.48	0.47
2:H:133:ASP:O	2:H:135:ARG:NE	2.44	0.47
1:D:176:TRP:O	1:D:180:ILE:HG12	2.15	0.47
2:H:216:LYS:HA	2:H:216:LYS:HD3	1.65	0.47
1:C:318:VAL:HG22	1:C:334:ARG:HD3	1.97	0.47
2:F:81:PHE:CZ	2:F:223:VAL:HG11	2.50	0.47
1:A:6:LEU:HD22	1:A:58:HIS:CE1	2.50	0.47
1:A:79:PRO:HB2	1:A:82:ILE:HD13	1.97	0.47
2:E:144:ASN:OD1	2:E:144:ASN:N	2.48	0.47
1:A:218:SER:OG	1:A:311:GLU:OE2	2.24	0.47
1:A:238:GLN:OE1	1:A:289:TRP:NE1	2.39	0.47
1:D:402:SER:OG	2:E:159:GLU:OE2	2.27	0.47
2:G:195:PRO:HG2	2:G:218:MET:HE1	1.96	0.47
1:C:358:GLN:NE2	2:G:39:GLU:OE2	2.47	0.47
1:B:418:GLU:OE2	2:H:68:THR:OG1	2.32	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:197:LYS:HG3	2:E:221:GLU:HB3	1.95	0.46
1:A:290:LYS:NZ	1:A:402:SER:O	2.37	0.46
2:E:15:VAL:HG22	2:E:25:PHE:HB3	1.97	0.46
1:B:28:TYR:CZ	1:B:49:VAL:HG21	2.50	0.46
2:G:179:LYS:HG2	2:G:213:PHE:CE1	2.50	0.46
2:H:70:ILE:HD13	2:H:73:GLN:HE22	1.81	0.46
1:D:196:ASN:OD1	1:D:199:ASN:HB3	2.15	0.46
2:G:201:THR:OG1	2:G:202:GLY:N	2.49	0.46
2:F:87:ILE:HG22	2:F:199:ILE:HD12	1.96	0.46
2:G:150:SER:HB2	2:G:180:THR:HG23	1.96	0.46
1:C:44:TYR:CD2	1:C:251:LYS:HG2	2.51	0.46
1:A:243:ASP:OD1	1:A:247:ARG:NH2	2.48	0.46
1:A:269:VAL:HG23	1:A:270:VAL:HG22	1.97	0.46
1:B:221:ARG:NH2	1:B:305:GLU:OE2	2.38	0.46
1:C:199:ASN:OD1	1:C:200:LEU:N	2.49	0.46
1:B:24:VAL:HG13	1:B:25:PRO:HD3	1.96	0.46
1:D:280:THR:OG1	1:D:306:GLU:OE1	2.31	0.45
2:H:84:VAL:HG12	2:H:85:ASP:H	1.81	0.45
2:G:196:ARG:O	2:G:220:VAL:HG11	2.17	0.45
1:D:137:VAL:HB	1:D:159:VAL:HG22	1.98	0.45
2:H:54:THR:OG1	2:H:69:GLU:HB3	2.16	0.45
1:B:221:ARG:HH11	1:B:233:VAL:HG11	1.82	0.45
1:C:283:PRO:HB2	1:C:309:ILE:HG13	1.99	0.45
1:D:408:GLN:O	1:D:412:ARG:HD3	2.16	0.45
1:D:53:HIS:CD2	1:D:129:PRO:HD2	2.51	0.45
2:H:52:VAL:HG23	2:H:240:ALA:HB2	1.98	0.45
1:D:75:CYS:HA	3:D:501:BJ8:S5	2.57	0.45
1:B:169:GLU:O	1:B:173:ARG:HG3	2.17	0.45
1:A:412:ARG:NH2	2:G:112:ASN:O	2.50	0.45
1:B:28:TYR:CE1	1:B:49:VAL:HG21	2.51	0.45
1:C:362:VAL:HG12	1:C:364:GLY:H	1.83	0.44
1:B:231:LEU:HD11	1:B:289:TRP:HB3	1.98	0.44
1:D:100:THR:HB	1:D:121:GLY:HA3	1.99	0.44
2:F:90:LEU:HD13	2:F:95:ALA:HB2	1.99	0.44
1:A:123:LYS:O	1:A:123:LYS:NZ	2.35	0.44
2:E:215:GLN:OE1	2:E:222:VAL:N	2.51	0.44
2:G:211:ARG:HD2	2:G:224:VAL:HG21	1.99	0.44
1:D:73:THR:HG21	1:D:78:VAL:HG21	1.99	0.44
1:B:221:ARG:HH21	1:B:303:VAL:C	2.25	0.44
1:C:233:VAL:HG12	1:C:237:ASN:HD21	1.83	0.44
2:G:101:ASP:OD2	2:G:105:ARG:NE	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:96:GLY:HA2	1:C:401:PHE:CE2	2.53	0.44
2:F:84:VAL:O	2:F:100:LYS:HE2	2.18	0.44
2:G:130:ILE:HG22	2:G:132:ASP:H	1.83	0.44
2:H:110:LEU:HD13	2:H:192:LEU:HD21	2.00	0.44
1:C:105:GLU:HG2	1:C:117:LYS:HD2	1.99	0.44
1:C:344:THR:HA	1:C:345:PRO:HA	1.86	0.44
2:H:51:LEU:HD23	2:H:51:LEU:H	1.83	0.44
1:A:283:PRO:HB2	1:A:309:ILE:HG13	1.99	0.43
1:B:64:LYS:NZ	1:B:135:ASP:OD1	2.45	0.43
1:C:395:LEU:HD12	1:C:416:PHE:HB2	2.00	0.43
1:C:284:GLN:HE22	1:C:292:HIS:HE1	1.66	0.43
2:F:233:ALA:O	2:F:236:ALA:HB3	2.19	0.43
1:B:56:ARG:HD2	1:B:56:ARG:HA	1.82	0.43
1:B:61:TYR:O	1:B:65:GLN:HG2	2.18	0.43
2:H:163:LEU:O	2:H:168:THR:OG1	2.26	0.43
2:H:233:ALA:O	2:H:236:ALA:HB3	2.18	0.43
1:C:367:HIS:CE1	1:C:380:SER:HB3	2.53	0.43
2:H:5:LEU:HD21	2:H:7:LEU:HG	1.99	0.43
1:C:408:GLN:O	1:C:412:ARG:HD3	2.19	0.43
1:D:317:LEU:C	1:D:334:ARG:HG3	2.43	0.43
1:A:56:ARG:NH1	1:A:59:GLU:OE1	2.49	0.43
1:A:276:ARG:HE	1:A:303:VAL:HG11	1.84	0.43
2:G:31:ARG:C	2:G:33:GLU:H	2.27	0.43
2:F:146:VAL:O	2:F:176:GLY:HA3	2.18	0.43
1:C:94:CYS:O	1:C:285:PRO:HB3	2.19	0.43
2:H:86:TYR:O	2:H:199:ILE:HD12	2.19	0.43
2:E:183:LYS:HB2	2:E:183:LYS:HE2	1.90	0.43
2:F:74:ALA:HB2	2:F:98:ILE:HD12	2.01	0.43
1:A:157:ILE:HG13	1:A:158:PRO:HD2	2.01	0.43
1:B:412:ARG:NH1	2:H:111:MET:HB3	2.34	0.43
1:A:340:CYS:SG	1:A:341:ALA:N	2.92	0.42
1:B:275:PRO:HD2	1:B:300:GLY:HA2	2.00	0.42
1:D:217:LEU:HD21	1:D:233:VAL:HG13	2.01	0.42
2:F:121:ARG:O	2:F:121:ARG:HD3	2.19	0.42
1:A:100:THR:HG21	1:A:122:PHE:CD1	2.54	0.42
2:F:30:THR:HG22	2:F:32:TYR:H	1.83	0.42
1:B:32:PHE:O	1:B:38:ARG:NE	2.50	0.42
1:B:33:LEU:HA	1:B:38:ARG:HH21	1.84	0.42
1:C:105:GLU:OE2	1:C:117:LYS:NZ	2.36	0.42
2:H:15:VAL:HG23	2:H:230:VAL:HG12	2.00	0.42
1:D:35:ARG:HA	1:D:36:PRO:HD3	1.91	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:ASP:O	1:B:49:VAL:HG22	2.19	0.42
1:B:216:ARG:NH1	1:B:257:ASP:OD1	2.46	0.42
2:E:183:LYS:NZ	2:E:217:GLU:HB3	2.35	0.42
2:E:136:ASP:HB2	2:E:209:ALA:HB2	2.02	0.42
1:A:216:ARG:O	1:A:220:LEU:HD23	2.20	0.42
1:B:397:LEU:HD23	1:B:397:LEU:HA	1.91	0.42
2:H:150:SER:OG	2:H:184:ARG:HD3	2.19	0.42
2:F:203:GLY:O	2:F:206:LYS:HG2	2.20	0.42
2:H:142:GLU:OE2	2:H:171:ARG:NE	2.53	0.42
2:H:28:VAL:HG12	2:H:41:LEU:HG	2.02	0.41
1:A:166:GLN:HG2	1:B:147:LYS:HZ3	1.85	0.41
1:A:344:THR:HA	1:A:345:PRO:HA	1.87	0.41
1:D:412:ARG:HD2	2:F:111:MET:SD	2.59	0.41
2:E:174:ILE:HD13	2:E:174:ILE:HA	1.94	0.41
2:F:227:GLU:HA	2:F:228:PRO:HD3	1.79	0.41
1:A:350:LEU:HD12	1:A:350:LEU:HA	1.91	0.41
1:A:155:GLU:OE2	1:B:385:ARG:NH1	2.53	0.41
2:H:237:ALA:O	2:H:240:ALA:HB3	2.21	0.41
2:E:16:ILE:O	2:E:23:TYR:HA	2.21	0.41
1:A:340:CYS:HB3	1:A:343:PHE:HE2	1.84	0.41
1:A:164:LEU:HD12	1:A:165:PRO:HD2	2.01	0.41
2:G:49:ARG:HD2	2:G:49:ARG:HA	1.93	0.41
1:C:395:LEU:HD22	1:C:397:LEU:HG	2.01	0.41
1:D:221:ARG:HH12	1:D:304:GLY:HA2	1.86	0.41
2:E:1:MET:N	2:E:17:ILE:O	2.41	0.41
2:G:143:GLU:OE1	2:G:143:GLU:N	2.53	0.41
1:C:211:ARG:NH2	1:C:313:TYR:O	2.52	0.41
2:E:5:LEU:HB3	2:E:53:VAL:HG23	2.03	0.41
2:E:198:LEU:HD22	2:E:218:MET:HE2	2.02	0.41
2:H:101:ASP:OD1	2:H:105:ARG:HG3	2.21	0.41
1:C:61:TYR:O	1:C:65:GLN:HG2	2.21	0.41
2:F:10:THR:C	2:F:30:THR:HB	2.46	0.41
1:A:16:LYS:HD2	1:A:17:HIS:N	2.36	0.41
1:B:15:GLU:OE1	1:B:15:GLU:N	2.30	0.41
2:G:85:ASP:HB3	2:G:100:LYS:O	2.21	0.41
2:H:13:LYS:HE2	2:H:229:SER:HB3	2.02	0.41
2:H:65:LYS:HG2	2:H:67:VAL:HG23	2.03	0.41
2:H:81:PHE:CE2	2:H:223:VAL:HG11	2.56	0.41
2:F:207:TYR:HA	2:F:208:PRO:HD3	1.90	0.41
1:B:21:LEU:HD23	1:B:21:LEU:HA	1.92	0.41
1:B:190:LYS:HD2	1:B:190:LYS:HA	1.81	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:43:LYS:HD3	2:G:43:LYS:HA	1.94	0.41
1:C:73:THR:HG22	1:C:138:VAL:HB	2.02	0.40
1:D:341:ALA:O	1:D:375:PRO:HB3	2.21	0.40
1:A:278:LEU:HB2	1:A:362:VAL:HG21	2.03	0.40
1:B:228:ILE:HD12	1:B:259:LEU:HD11	2.03	0.40
2:G:137:GLU:CD	2:G:137:GLU:H	2.28	0.40
2:H:13:LYS:HB3	2:H:230:VAL:HG13	2.03	0.40
2:H:136:ASP:OD2	2:H:137:GLU:HG3	2.21	0.40
1:A:6:LEU:HD22	1:A:58:HIS:ND1	2.35	0.40
1:C:319:GLU:C	1:C:330:ASN:HD22	2.28	0.40
2:E:133:ASP:OD2	2:E:133:ASP:N	2.53	0.40
1:C:10:LEU:HD23	1:C:10:LEU:HA	1.89	0.40
1:D:249:ALA:O	1:D:252:VAL:HG12	2.22	0.40
1:D:350:LEU:HD12	1:D:350:LEU:HA	1.92	0.40
2:G:149:ASN:OD1	2:G:150:SER:N	2.55	0.40
2:H:134:TYR:CE2	2:H:138:ASP:HB2	2.57	0.40
2:H:201:THR:OG1	2:H:202:GLY:N	2.54	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	420/422 (100%)	397 (94%)	23 (6%)	0	100	100
1	B	418/422 (99%)	382 (91%)	36 (9%)	0	100	100
1	C	420/422 (100%)	390 (93%)	27 (6%)	3 (1%)	18	48
1	D	421/422 (100%)	389 (92%)	32 (8%)	0	100	100
2	E	241/243 (99%)	225 (93%)	16 (7%)	0	100	100
2	F	239/243 (98%)	210 (88%)	27 (11%)	2 (1%)	16	44
2	G	241/243 (99%)	225 (93%)	16 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	239/243 (98%)	202 (84%)	37 (16%)	0	100	100
All	All	2639/2660 (99%)	2420 (92%)	214 (8%)	5 (0%)	43	70

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	34	PRO
1	C	285	PRO
2	F	228	PRO
1	C	240	ALA
1	C	287	PRO

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	364/363 (100%)	364 (100%)	0	100	100
1	B	362/363 (100%)	362 (100%)	0	100	100
1	C	364/363 (100%)	364 (100%)	0	100	100
1	D	365/363 (101%)	365 (100%)	0	100	100
2	E	196/196 (100%)	196 (100%)	0	100	100
2	F	194/196 (99%)	194 (100%)	0	100	100
2	G	196/196 (100%)	196 (100%)	0	100	100
2	H	194/196 (99%)	194 (100%)	0	100	100
All	All	2235/2236 (100%)	2235 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	237	ASN

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Mol	Chain	Res	Type
1	C	292	HIS
1	C	346	ASN
1	D	11	ASN
1	D	199	ASN
1	D	237	ASN
1	D	253	ASN
2	F	78	HIS
2	F	215	GLN
1	A	65	GLN
1	A	237	ASN
1	A	408	GLN
1	B	3	ASN
1	B	58	HIS
1	B	98	GLN
1	B	154	ASN
1	B	166	GLN
1	B	237	ASN
1	B	284	GLN
1	B	292	HIS
2	G	144	ASN
2	G	215	GLN
2	H	62	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SF4	E	301	-	0,12,12	-	-	-		
4	SF4	G	301	-	0,12,12	-	-	-		
3	BJ8	B	501	-	6,26,26	3.92	4 (66%)	-		
5	SO4	F	301	-	4,4,4	0.23	0	6,6,6	0.08	0
5	SO4	E	302	-	4,4,4	0.25	0	6,6,6	0.09	0
3	BJ8	D	501	-	6,26,26	3.42	5 (83%)	-		
5	SO4	G	302	-	4,4,4	0.23	0	6,6,6	0.07	0
3	BJ8	C	501	-	6,26,26	5.44	5 (83%)	-		
3	BJ8	A	501	-	6,26,26	4.67	4 (66%)	-		

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	SF4	E	301	-	-	-	0/6/5/5
4	SF4	G	301	-	-	-	0/6/5/5
3	BJ8	B	501	-	-	-	0/12/10/10
3	BJ8	D	501	-	-	-	0/12/10/10
3	BJ8	C	501	-	-	-	0/12/10/10
3	BJ8	A	501	-	-	-	0/12/10/10

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	501	BJ8	S5-FE8	10.18	2.44	2.28
3	A	501	BJ8	S5-FE8	9.61	2.43	2.28
3	B	501	BJ8	S5-FE8	7.62	2.40	2.28
3	C	501	BJ8	S2-FE1	6.04	2.37	2.28
3	D	501	BJ8	S2-FE1	5.29	2.36	2.28
3	B	501	BJ8	S2-FE1	4.38	2.35	2.28
3	C	501	BJ8	S4-FE1	-4.36	2.21	2.28
3	D	501	BJ8	S7-FE8	4.09	2.34	2.28
3	A	501	BJ8	S6-FE8	3.92	2.34	2.28
3	D	501	BJ8	S5-FE8	3.77	2.34	2.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	BJ8	S7-FE8	-3.45	2.22	2.28
3	C	501	BJ8	S6-FE8	3.08	2.32	2.28
3	C	501	BJ8	S7-FE8	-2.87	2.23	2.28
3	B	501	BJ8	S6-FE8	2.87	2.32	2.28
3	A	501	BJ8	S3-FE1	-2.39	2.24	2.28
3	B	501	BJ8	S4-FE1	-2.10	2.24	2.28
3	D	501	BJ8	S4-FE1	-2.05	2.24	2.28
3	D	501	BJ8	S6-FE8	2.00	2.31	2.28

There are no bond angle outliers.

There are no chirality outliers.

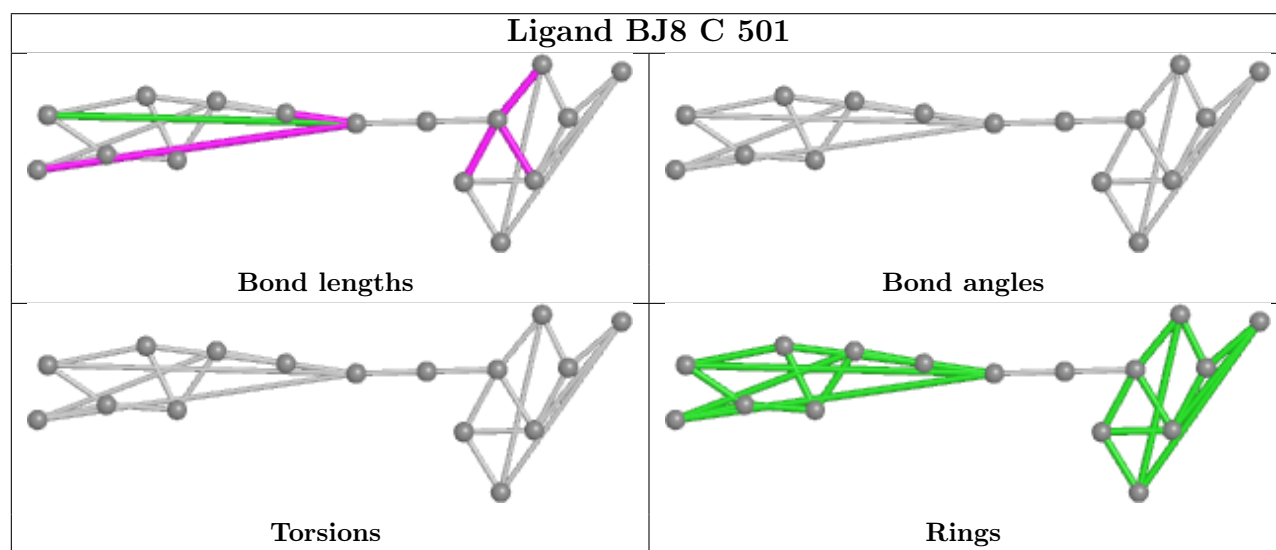
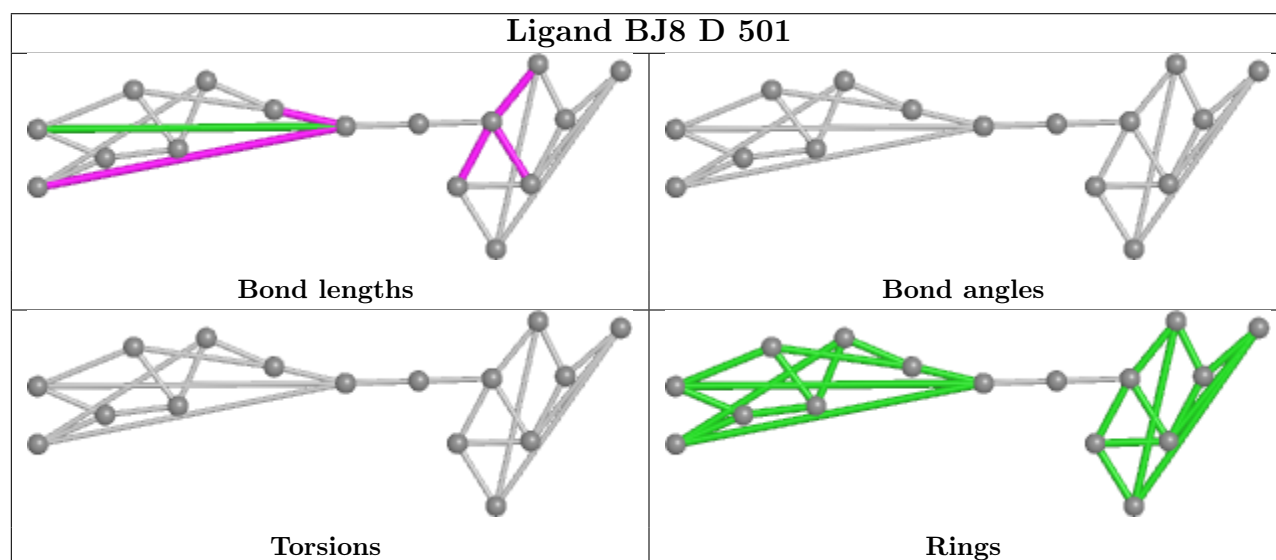
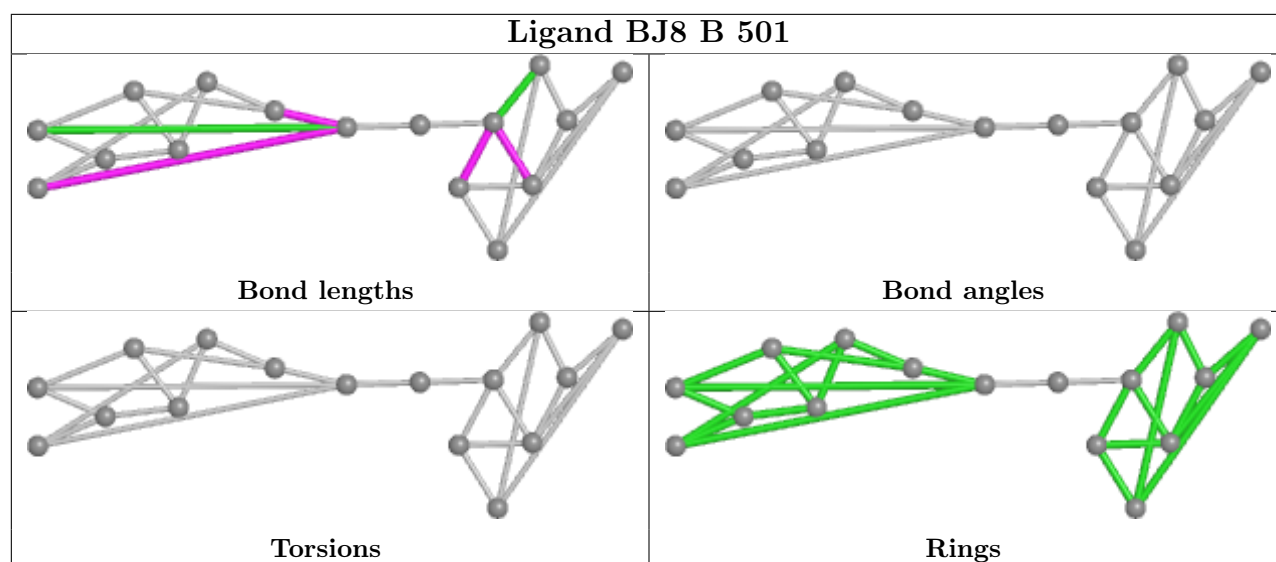
There are no torsion outliers.

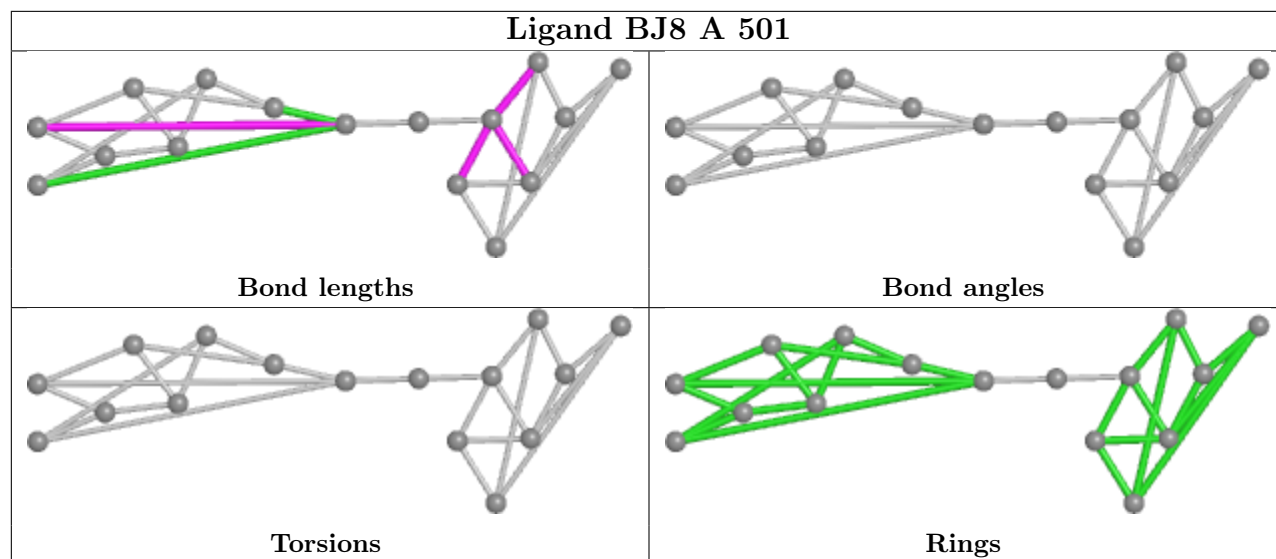
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	302	SO4	1	0
3	D	501	BJ8	1	0
5	G	302	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	421/422 (99%)	0.28	7 (1%) 69 50	26, 42, 62, 83	2 (0%)
1	B	419/422 (99%)	0.46	10 (2%) 59 40	32, 55, 71, 89	1 (0%)
1	C	421/422 (99%)	0.24	5 (1%) 76 58	24, 39, 55, 65	2 (0%)
1	D	420/422 (99%)	0.41	12 (2%) 53 36	27, 52, 72, 94	3 (0%)
2	E	243/243 (100%)	0.43	7 (2%) 53 36	32, 56, 98, 115	5 (2%)
2	F	241/243 (99%)	1.26	46 (19%) 3 3	35, 93, 125, 133	6 (2%)
2	G	243/243 (100%)	0.66	11 (4%) 38 25	32, 61, 101, 125	5 (2%)
2	H	241/243 (99%)	1.32	50 (20%) 2 2	61, 99, 134, 145	5 (2%)
All	All	2649/2660 (99%)	0.55	148 (5%) 30 20	24, 53, 113, 145	29 (1%)

All (148) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	139	LEU	5.1
2	H	232	ALA	5.1
2	H	229	SER	4.6
2	F	133	ASP	4.5
2	H	118	GLY	3.9
2	H	15	VAL	3.8
1	D	77	TYR	3.8
1	A	16	LYS	3.7
2	F	209	ALA	3.7
2	H	84	VAL	3.7
2	H	115	CYS	3.7
1	A	70	VAL	3.6
2	F	189	ALA	3.6
2	F	222	VAL	3.6
1	D	281	GLY	3.5
2	E	21	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
2	H	214	LEU	3.4
2	F	234	LEU	3.4
2	F	198	LEU	3.4
2	E	236	ALA	3.4
2	H	218	MET	3.4
2	H	117	ALA	3.4
2	F	51	LEU	3.3
2	F	214	LEU	3.3
2	H	180	THR	3.2
2	H	224	VAL	3.2
2	H	210	LEU	3.1
2	G	138	ASP	3.1
2	F	155	PHE	3.1
2	H	77	CYS	3.1
2	H	122	PHE	3.1
2	E	174	ILE	3.1
1	D	136	TYR	3.0
2	F	138	ASP	3.0
1	B	93	LEU	3.0
1	B	193	VAL	3.0
2	H	53	VAL	3.0
2	F	70	ILE	3.0
1	D	195	LEU	3.0
2	F	239	ILE	3.0
2	F	115	CYS	2.9
2	H	24	THR	2.9
2	H	51	LEU	2.9
1	C	373	CYS	2.9
2	F	141	ASN	2.9
2	H	16	ILE	2.9
2	H	41	LEU	2.9
2	H	234	LEU	2.9
2	F	204	GLY	2.9
2	F	3	ALA	2.8
1	B	418	GLU	2.8
2	F	74	ALA	2.8
2	H	100	LYS	2.8
2	G	173	VAL	2.8
2	H	40	LEU	2.8
2	F	50	ASN	2.7
2	H	134	TYR	2.7
2	H	61	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
2	H	219	GLY	2.7
2	H	165	ALA	2.7
2	G	118	GLY	2.6
1	D	184	ALA	2.6
2	H	3	ALA	2.6
2	F	225	PRO	2.6
2	G	40	LEU	2.6
2	F	118	GLY	2.6
2	H	152	CYS	2.5
2	E	139	LEU	2.5
1	A	126	ARG	2.5
2	F	106	VAL	2.5
2	F	107	VAL	2.5
2	E	155	PHE	2.5
2	F	23	TYR	2.5
2	H	158	SER	2.5
2	F	213	PHE	2.5
2	F	4	GLY	2.5
2	H	127	LEU	2.5
2	H	32	TYR	2.5
2	F	154	VAL	2.5
1	A	6	LEU	2.4
2	F	195	PRO	2.4
2	H	230	VAL	2.4
2	F	87	ILE	2.4
1	D	142	THR	2.4
2	H	195	PRO	2.4
2	F	219	GLY	2.4
2	F	98	ILE	2.4
1	A	1	MET	2.4
2	H	213	PHE	2.4
2	H	28	VAL	2.3
2	H	35	VAL	2.3
2	F	6	ASP	2.3
1	A	15	GLU	2.3
1	B	209	LYS	2.3
1	D	114	PRO	2.3
2	H	139	LEU	2.3
1	B	84	ASN	2.3
2	G	238	LEU	2.3
2	H	231	THR	2.3
1	C	75	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
2	H	209	ALA	2.2
2	E	134	TYR	2.2
2	F	158	SER	2.2
1	A	37	ASN	2.2
1	D	259	LEU	2.2
2	F	85	ASP	2.2
2	H	189	ALA	2.2
1	B	138	VAL	2.2
1	C	1	MET	2.2
1	D	94	CYS	2.2
1	B	321	ALA	2.2
2	F	77	CYS	2.2
2	F	134	TYR	2.2
2	H	73	GLN	2.2
2	G	156	ALA	2.2
2	F	231	THR	2.1
2	G	220	VAL	2.1
2	F	218	MET	2.1
2	F	190	GLU	2.1
1	D	192	GLY	2.1
2	G	231	THR	2.1
1	D	204	ILE	2.1
2	H	182	ALA	2.1
2	G	158	SER	2.1
2	G	155	PHE	2.1
2	H	52	VAL	2.1
1	B	204	ILE	2.1
2	F	68	THR	2.1
2	H	60	ALA	2.1
2	H	225	PRO	2.1
2	H	220	VAL	2.1
2	F	99	LYS	2.1
1	B	308	CYS	2.1
2	F	54	THR	2.1
2	H	110	LEU	2.0
2	H	212	LEU	2.0
2	F	117	ALA	2.0
2	G	165	ALA	2.0
2	F	81	PHE	2.0
2	F	53	VAL	2.0
2	E	161	ILE	2.0
1	D	3	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
2	F	233	ALA	2.0
1	C	77	TYR	2.0
1	C	405	ASP	2.0
1	B	75	CYS	2.0
2	H	241	ARG	2.0
2	H	148	ILE	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

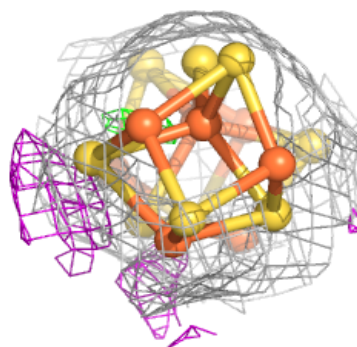
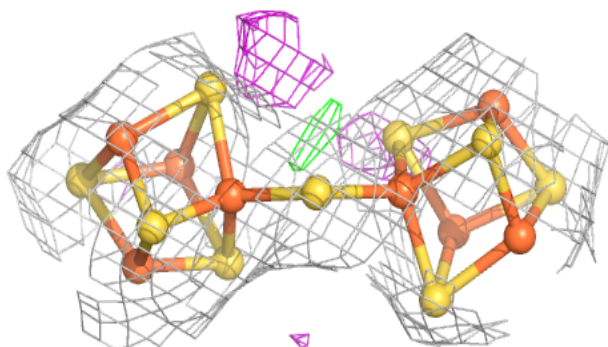
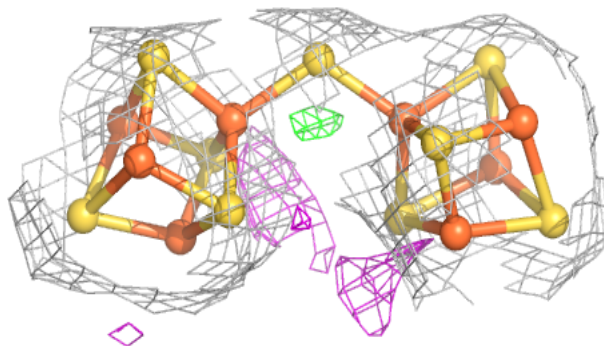
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SO4	G	302	5/5	0.80	0.15	61,67,69,73	0
5	SO4	F	301	5/5	0.82	0.12	88,89,90,104	0
5	SO4	E	302	5/5	0.91	0.11	66,66,77,79	0
4	SF4	E	301	8/8	0.96	0.10	66,80,94,98	0
4	SF4	G	301	8/8	0.97	0.08	62,76,82,93	0
3	BJ8	A	501	17/17	0.97	0.07	24,34,48,50	0
3	BJ8	C	501	17/17	0.98	0.05	23,33,41,45	0
3	BJ8	D	501	17/17	0.98	0.05	21,34,54,57	0
3	BJ8	B	501	17/17	0.99	0.06	25,34,48,48	0

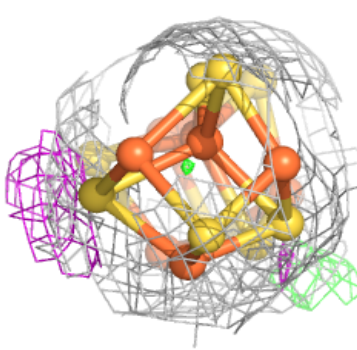
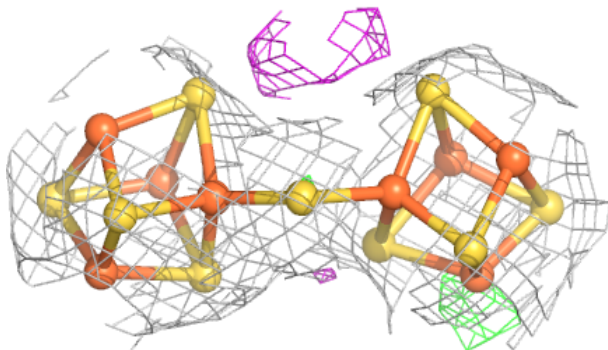
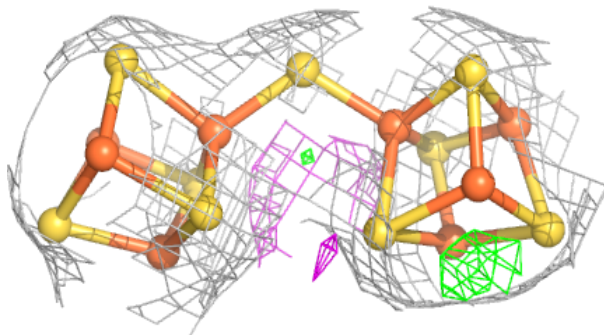
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around BJ8 A 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

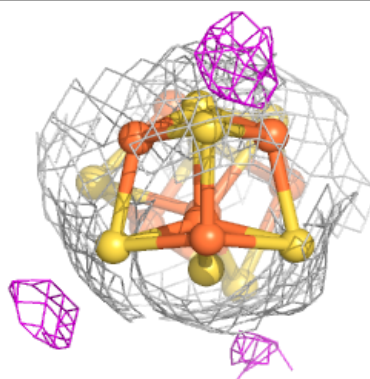
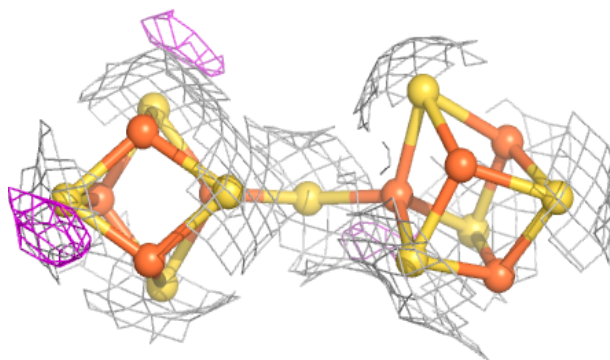
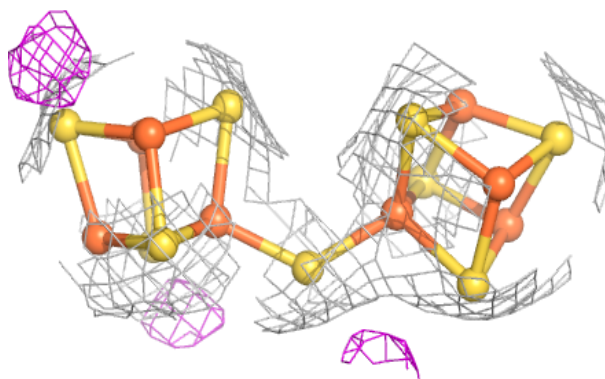
**Electron density around BJ8 C 501:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

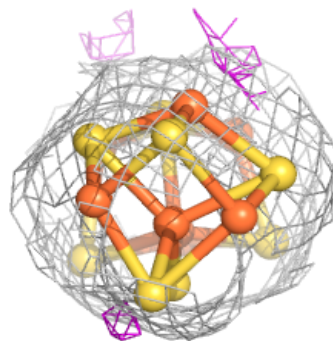
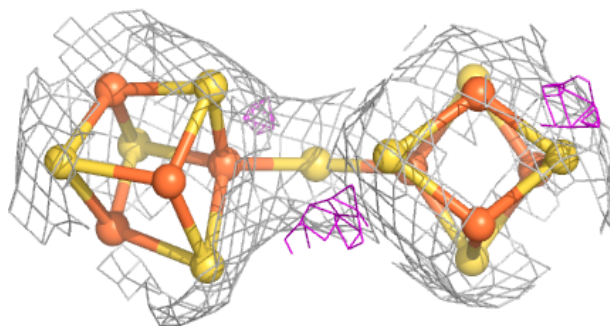
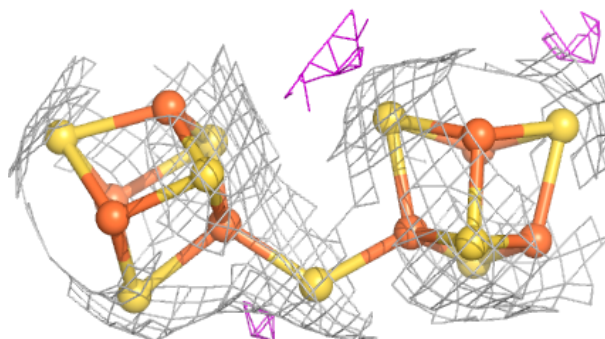


Electron density around BJ8 D 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around BJ8 B 501:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.